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**QUARTERLY MONITORING REPORT
ACTIVE TREATMENT SYSTEMS
FOURTH QUARTER 2006**

**AMERICAN CHEMICAL SERVICE NPL SITE
GRIFFITH, INDIANA**

MWH File No. 4050577

Prepared For:

**American Chemical Service NPL Site RD/RA Executive Committee
Griffith, Indiana**

Prepared By:

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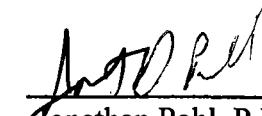
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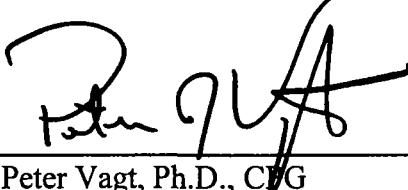


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ACRONYMS AND ABBREVIATIONS

AS	Air Sparge
AMSL	Above Mean Sea Level
BOD	Biological Oxygen Demand
BW	Barrier Wall
BWES	Barrier Wall Extraction System
cfm	cubic feet per minute
DL	Detection Limit
DPE	Dual Phase Extraction
GAC	Granular Activated Carbon
Global	Global Technologies
GWTP	Groundwater Treatment Plant
"Hg	Inches of mercury
"H ₂ O	Inches of water
IDEM	Indiana Department of Environmental Management
K-P	Kapica Pazmey
lb/hr	Pounds per hour
LDC	Laboratory Data Consultants
mg/kg	Milligrams per kilogram
mg/L	Milligrams per liter
NC	Not Calculated
ND	Not Detected
NE	No Effluent Limit Established
NS	Not Sampled
OFCA	Off-Site Containment Area
PCBs	Polychlorinated Biphenyls
ppm	Parts per million
PGCS	Perimeter Groundwater Containment System
PSVP	Performance Standard Verification Plan
QAPP	Quality Assurance Project Plan
QA/QC	Quality Assurance/Quality Control
SBPA	Still Bottoms Pond Area
SVOC	Semi-Volatile Organic Compounds
T-102	Aeration Equalization Tank (Tank – 102)
TOC	Top of Casing
TOIC	Top of Inner Casing
TOSG	Top of Staff Gauge
TSS	Total Suspended Solids
µg	Micrograms
µg/L	Micrograms per liter
U.S. EPA	United States Environmental Protection Agency
VOCs	Volatile Organic Compounds

1.0 INTRODUCTION

MWH Americas, Inc. (MWH), on behalf of the American Chemical Service (ACS) Remedial Design/Remedial Action (RD/RA) Executive Committee, started up the on-site groundwater treatment system at the ACS National Priorities List (NPL) Site (ACS Site) in Griffith, Indiana on March 13, 1997. The groundwater treatment plant (GWTP) system was designed to treat groundwater from the Perimeter Groundwater Containment System (PGCS) and the Barrier Wall Extraction System (BWES). The original treatment consisted of a phase-separator for oil and free product removal, equalization tanks, an UV oxidation unit for destruction of organic constituents, and an air stripper to remove methylene chloride and other organics. The treatment also included a chemical precipitation and clarification unit to remove metals, a sand filter to remove suspended solids, and activated carbon vessels for final polishing of the treated groundwater before it was released to the west of the Site.

In 2001, an activated sludge treatment unit was added to the process to reduce the volatile and semivolatile organic compounds (VOCs and SVOCs) in the collected groundwater. The activated sludge treatment process also reduces the amount of activated carbon required to treat the water. An aerated equalization tank was also added to the GWTP in 2001 to remove VOCs from the collected groundwater, oxidize metals to increase metals removal efficiency in the chemical precipitation unit, and equalize groundwater flow through the GWTP. The activated sludge system and aeration tank have been fully integrated into the process along with the other upgrade components. Startup and optimization of the catalytic oxidizer/scrubber air treatment unit was also conducted during 2001.

The treated effluent from the treatment system is discharged to the nearby wetlands, west of the treatment system, in accordance with Agency approvals.

Operation of the In-situ Soil Vapor Extraction (ISVE) system for the Off-Site Containment Area (OFCA) and the Kapica-Pazmey (K-P) Area began on May 1, 2002. Operation of the ISVE system for the Still Bottoms Pond Area (SBPA) began in July 2003. The ISVE systems were designed to remove volatile and semi-volatile compounds from the subsurface media.

The Off-Site Area ISVE system consists of 42 ISVE wells, 3 air sparge wells, ISVE and air sparge blower systems, a thermal oxidizer/scrubber unit, and the associated mechanical and electrical components. Protocols and goals for the phased startup of the Off-Site System as defined in the Final Remedy (Montgomery Watson, 1999) were followed. In 2004, an additional blower unit was added to the Off-Site Area ISVE system to more effectively meet the design objectives of the system. The additional blower increased the capacity of the Off-Site ISVE system from 1000 to 2000 cubic feet per minute (cfm).

The SBPA ISVE system consists of 25 ISVE wells, 21 dual-phase extraction (DPE) wells, 6 air sparge wells, ISVE and air sparge blower systems, a thermal oxidizer/scrubber unit, and the associated mechanical and electrical components. During the first 12 months of system operation, the performance of the ISVE system was evaluated. Based on this evaluation, the

SBPA ISVE system was enhanced in accordance with the United States Environmental Protection Agency (U.S. EPA) and Indiana Department of Environmental Management (IDEM) approval by reconfiguring 18 of the ISVE wells to allow injection of air. Air for the injection wells is directed from blower ME-102/103 at the GWTP to the SBPA ISVE blower shed. The air injection system, which consists of three groups of five injection wells, began operation in December 2005. The air injection is scheduled to rotate among the three well groups on a monthly basis. Only one well group is operated at a time.

This report summarizes GWTP effluent analytical data and thermal oxidizer off-gas analytical data, ISVE process monitoring data, and water level gauging data collected from October 2006 through December 2006. The report also details modifications and upgrades that were made to the active treatment systems during the reporting period.

2.0 GWTP COMPLIANCE MONITORING

2.1 SAMPLING REQUIREMENTS

Effluent samples are collected on a regular schedule from the treatment system to demonstrate compliance with the discharge limits (Table 2.1) established by the Indiana Department of Environmental Management (IDEM) and the United States Environmental Protection Agency (U.S. EPA). The approved Performance Standard Verification Plan (PSVP) (Montgomery Watson, July 1997) requires quarterly effluent sampling for biochemical oxygen demand (BOD), total suspended solids (TSS), SVOCs, metals, and polychlorinated biphenyls (PCBs) in the system, and monthly effluent sampling for pH and VOCs, as tabulated below. In accordance with the PSVP, a full analysis effluent compliance sample was collected during October 2006 and analyzed for all of the analytes listed above. During November and December 2006, the monthly effluent compliance samples were analyzed for VOCs and pH only.

Sampling and analyses were performed in accordance with the approved Quality Assurance Project Plan (QAPP) (Montgomery Watson Harza, November 2001). Quality control measures were also instituted in accordance with the PSVP. The following table and paragraphs present details on sampling and analyses and also summarize the analytical data for the treatment system effluent.

Sampling Frequency Schedule – Groundwater Treatment System

Analytes	Cumulative Time From Startup*	Frequency
Flowrate	–	Continuous
BOD, TSS, SVOCs and Metals	181 days onward	Once per quarter
VOCs and pH	31 days onward	Once per month
PCBs	181 days onward	Once per quarter
PCBs in Sediment (one location)	–	Once per year

*Note: System operation began on March 13, 1997

2.2 EFFLUENT SAMPLING AND ANALYSES

Effluent samples were collected each month during the fourth quarter of 2006. Samples were collected on the following dates and analyzed for the listed analytes for this reporting period:

October 11, 2006	Full analysis (pH, TSS, BOD, Metals, VOCs, SVOCs, pentachlorophenol, and PCBs)
November 1, 2006	pH and VOCs
December 21, 2006	pH and VOCs

The above samples were collected directly from a sampling port on the effluent line of the treatment system. The samples were placed in contaminant-free containers, in accordance with the U.S. EPA Specifications and Guidance for Obtaining Contaminant-Free Sample Containers (U.S. EPA, 1992). Appropriate sample containers and preservatives, as specified in the QAPP, were used to collect and preserve the samples. Following sample collection, the temperature of the sample containers was maintained at or below 4° C in coolers. Chain-of-Custody forms were prepared to track the transfer of samples from the treatment system to the laboratories. In accordance with the approved QAPP, the effluent water samples were analyzed for the following parameters by the following analytical methods:

<u>Parameter</u>	<u>Analytical Method</u>
VOCs	SW-846 8260B
SVOCs	SW-846 8270C
Pentachlorophenol	SW-846 8270C and SIM
Pesticides/PCBs	EPA 608/SW-846 8081/8082
Metals (Excluding Mercury)	
General Water Quality	SW-846 6010
Parameters (TSS and BOD-5)	EPA 160.2 and 405.1
Mercury	SW-846 7470
pH	EPA 150.1

2.3 EFFLUENT ANALYTICAL RESULTS

2.3.1 GWTP Effluent Samples

The GWTP effluent monitoring data, summarized in Table 2.2, verify that the system effluent was compliant with the discharge limits summarized in Table 2.1. No effluent exceedences were reported in the October, November, or December samples.

Compuchem Laboratory of Cary, North Carolina performed the analysis of the samples. Laboratory Data Consultants (LDC) of Carlsbad, California performed third party data validation in accordance with the U.S. EPA National Functional Guidelines for Organic/Inorganic Data Review (U.S. EPA, February 1994 and October 1999). Validation qualifiers are listed in Table 2.2 and are written in the margin of the analytical data sheets provided in Appendix A.

2.3.2 Sediment Sample

Since 1998, MWH has collected an annual sediment sample and associated quality control samples from the GWTP outfall in accordance with the PSVP to help determine if PCB accumulation is occurring at the GWTP discharge location. The annual sediment sample for 2006 was collected on December 11th from the GWTP outfall location, shown on Figure 2.1. The sample was analyzed for PCBs by CompuChem and the data was validated by LDC.

The analytical data for the annual sediment samples for the past seven years are summarized in Table 2.3. Analytical data for the December 2006 sample are included in Appendix C. One aroclor, Aroclor-1248, was detected in the December 2006 sample. Although

concentrations remain below any action levels, results for this compound had been non-detect for previous sampling events. MWH will accelerate the schedule for collection of the 2007 sampling event to coincide with the next groundwater sampling event to be conducted in April.

3.0 ISVE SYSTEM MONITORING

3.1 THERMAL OXIDIZER OFF-GAS SAMPLING

During the fourth quarter of 2006, Thermal Oxidizer/Scrubber Unit 1 (Therm Ox 1) was used to treat vapors from the SBPA ISVE system and Thermal Oxidizer/Scrubber Unit 2 (Therm Ox 2) was used to treat vapors from the Off-Site ISVE system and T-102. VOC removal rates as well as the total VOCs removed are illustrated in Figure 3.1 and Figure 3.2, respectively. Compliance samples were collected from both thermal oxidizer/scrubber units on October 19th, November 2nd, and December 11th.

Influent and effluent off-gas samples were collected directly from sampling ports on the influent pipe to the thermal oxidizer and the discharge stack of the scrubber. One influent sample and one effluent sample were collected. A duplicate influent sample was also collected. The samples were collected to comply with the PSVP and QAPP and in accordance with laboratory guidelines. The VOC samples were collected using a Summa canister and the SVOC samples were collected in sorbent tubes.

Sampling Frequency Schedule – ISVE System

Startup	Weekly for a four week period
Post-Startup	Monthly in accordance with the IDEM Air Permit Equivalency

Following sample collection, the sorbent tubes were maintained at or below 4°C in coolers. Chain-of-Custody forms were prepared to track the transfer of samples from the treatment system to the laboratories for extraction and analysis. In accordance with the approved QAPP, the off-gas samples were analyzed by the following analytical methods:

<u>Parameter</u>	<u>Analytical Method</u>
VOCs	TO-14
SVOCs	TO-13

3.2 SAMPLING RESULTS

The influent and effluent off-gas data are collected to verify that the off-gas from both of the thermal oxidizers were less than the IDEM discharge limit of three pounds of VOCs per hour for October, November, and December. For example, the VOC discharge reported from the October 19, 2006 Therm Ox 1 sample was 0.015 pounds per hour, approximately 0.5 percent of the discharge limit. The VOC discharge from the October 19, 2006 Therm Ox 2 sample was 0.1429 pounds per hour, approximately five percent of the discharge limit. All of the results for November and December were within the same order of magnitude. Therefore, it can be concluded that the ISVE systems are performing well within discharge limits for air emissions. VOC discharge values for Therm Ox 1, Therm Ox 2, and the SBPA and Off-Site

ISVE system are presented in Tables 3.1 through 3.9. The analytical data sheets for the compliance samples are provided in Appendix B.

In addition to the off-gas data collected during the fourth quarter, MWH collected off-gas samples from the Off-Site ISVE system and the SBPA ISVE system influent lines. These samples were collected in order to comply with the PSVP.

Air Toxics Laboratories of Folsom, California analyzed the samples. The analytical results are summarized in Tables 3.1 through 3.18. MWH performed data validation in accordance with the QAPP and the National Functional Guidelines for Organic/Inorganic Data Review. Validation qualifiers are listed in the tables and are written in the margin of the analytical data sheets provided in Appendix B.

3.3 ISVE SYSTEM MONITORING

Performance monitoring of the ISVE system was conducted in accordance with the PSVP (Montgomery Watson, June 1999). Extracted vapor flow rates and vacuums at individual ISVE wells and headers were measured and recorded on a routine basis. Additionally, VOC concentrations were measured at individual wells and headers using a photoionization detector (PID).

The information collected during performance monitoring is used to evaluate and optimize the ISVE system. Data collected from the Off-Site ISVE system during the fourth quarter of 2006 are presented in Tables 3.19 and 3.20. Data that were collected from the SBPA ISVE system during the fourth quarter of 2006 are presented in Tables 3.21 and 3.22.

3.4 PRODUCT REMOVAL ACTIVITIES

Product removal activities were performed at two ISVE well locations in the SBPA during the fourth quarter 2006 (SVE-53 and SVE-72). A total of 38 gallons of liquid were removed from these wells. The product removal schedule for the fourth quarter is summarized in Table 3.23.

4.0 GWTP PROCESS MODIFICATIONS AND REPAIRS

4.1 GWTP PROCESS MODIFICATIONS

No modifications were made to the GWTP during the fourth quarter of 2006.

4.2 GWTP REPAIRS AND MAINTENANCE

The following maintenance was made to the GWTP during the fourth quarter of 2006:

- The temperature in the biotank continues to be maintained at approximately 58 degrees in order to sustain a viable microbial population. The heat exchanger used to warm the biotank influent water was inspected during the annual maintenance event in October and found to be intact with minimal noticeable corrosion.
- MWH is currently evaluating options to use the free product collected from the SBPA ISVE wells as a fuel source for system operations. By using the free product as a fuel source, MWH can reduce natural gas usage. Options include injecting the product into the thermal oxidizers or using it in an oil heater to supplement facility heating.
- Recent groundwater level monitoring results indicate that levels in both the On-Site and Off-Site Areas have elevated during the gauging event on November 22nd. MWH is evaluating the performance of extraction wells to ensure correct operation and will continue to monitor groundwater levels.

5.0 ISVE PROCESS MODIFICATIONS AND REPAIRS

5.1 ISVE PROCESS MODIFICATIONS

The following modifications were made to the SBPA ISVE system during the fourth quarter of 2006:

- Three sets of five air injection wells ran at the ACS site throughout the fourth quarter 2006. On October 19, 2006, MWH switched the air injection wells from Group 3 (SVE-44, SVE-59, SVE-77, SVE-80, and SVE-84) to Group 1 (SVE-50, SVE-54, SVE-73, SVE-79, and SVE-81). Group 1 operated until November 29, 2006 when MWH switched over to Group 2 (SVE-49, SVE-51, SVE-65, SVE-71, and SVE-82). On December 21, 2006, MWH was at the Site to switch the air injection wells from Group 2 to Group 3. MWH will continue to rotate among the three groups of air injection wells on a monthly basis.
- MWH is currently evaluating a flow meter for improved vapor flow monitoring from the ISVE systems and thermal oxidizers. MWH has ordered a FCI ST98 thermal mass flow meter that will be installed on one of the thermal oxidizer influent pipes. The meter is anticipated to be installed during the week of January 8, 2007. MWH will evaluate the meter's performance for approximately two to three weeks to determine if the meter is appropriate for the application. If performance is satisfactory, three additional flow meters will be procured and installed to track vapor flow through the entire system.

No modifications were made to the Off-Site ISVE system during the fourth quarter of 2006.

5.2 ISVE REPAIRS AND MAINTENANCE

The following repairs were made to the ISVE system during the fourth quarter of 2006:

- Annual maintenance of the air sparge compressors was performed during November.
- Thermal Oxidizer 1 (ThermOx 1) was shut down during the month of November for routine maintenance activities.
- Thermal Oxidizer 2 (ThermOx 2) was shut down during November due to a malfunctioning flow switch on the scrubber recirculation line. The spare switch stored at the Site failed to operate correctly when it was installed, so MWH ordered a replacement switch as well as a new spare switch to be stored at the Site.

6.0 PGCS AND BWES GAUGING ACTIVITIES

During the operational time frame of the GWTP in the fourth quarter of 2006, the PGCS groundwater extraction trenches were operated in "auto" mode. In "auto" mode, the PGCS extraction wells pump continuously unless there is a low water level in individual extraction wells or a high water level in the Aeration Equalization Tank (T-102). This mode is used to control the flowrate through the treatment system, while at the same time creating an inward gradient along the PGCS trench. The GWTP also received influent from the On-Site and Off-Site components of the BWES, the SBPA DPE wells, and MW-56 during the fourth quarter of 2006. The pump in MW-10C malfunctioned. Therefore, pumping did not occur at this location during the fourth quarter 2006. The pump for MW-10C will be brought back online upon completion of the Lower Aquifer Pumping System.

In accordance with the PSVP, a discussion on the effect of the PGCS and BWES on the water table near the Site is presented in each quarterly monitoring report. This section summarizes the groundwater elevations at the Site during October, November, and December 2006. Groundwater elevation measurements were collected throughout the Site on December 22, 2006 as part of the groundwater monitoring program. The groundwater elevations are listed in Table 6.1 and the resulting water table contours outside the barrier wall are shown on Figure 6.1.

The barrier wall was constructed to contain the contaminated zone under the Site and the BWES was installed to extract groundwater from within the barrier wall and dewater the Site for the ISVE system. Eight pairs of piezometers were installed, with one piezometer of each pair on either side of the barrier wall, spaced along the barrier wall alignment. This allows measurement and tracking of water levels in order to document that the barrier wall is serving its designed function.

Table 6.1, BWES Water Level and Piezometer Pairs, presents the groundwater elevations inside and outside the barrier wall on December 22, 2006. The groundwater elevations are plotted on Figure 6.2. The groundwater elevation measurements inside the barrier wall range from 3.09 to 11.48 feet lower than levels outside the barrier wall. In general, the data demonstrates that the barrier wall is successfully performing the intended function of isolating and protecting the groundwater outside the barrier wall from the source areas of the Site inside the barrier wall. MWH will continue to collect water level measurements quarterly across the Site as required in the PSVP.

As part of the optimization of the GWTP and BWES upgrades, MWH began active dewatering of the Off-Site Area through increased groundwater pumping rates on September 25, 2001. Active dewatering of the SBPA (on-site area) began on February 11, 2003 with the addition of the DPE wells. Water levels were measured throughout the quarter at piezometer locations (P29, P31, P32, P36, and P49) in the On-Site Area and at piezometers (P96, P110, P112, P113, P114, P116, P118) and three air sparge (AS) wells (AS-7, AS-8, and AS-9) in the Off-Site Area. These locations are shown on Figure 6.3. The water level trend data from these piezometers and AS wells for the fourth

quarter 2006 are depicted graphically on Figures 6.4 and 6.5, which also show the target water elevations for each area. In the SBPA, the target water level is 629 feet amsl. Water levels in two piezometer locations (P-29 and P-32) have been drawn down to below the bottom of the screens in these wells throughout the fourth quarter 2006. Therefore, our depth to water measurements at these locations show straight-line measurements of the bottom of the wells. The other three locations had water levels that were drawn down to the below the well screen for the first half of the fourth quarter. However, during the second half of the quarter, water levels varied from approximately 625 feet amsl to 632 feet amsl. This represents an increase in the average water levels from the third quarter 2006.

In the Off-Site ISVE area, the target water level is 626 feet amsl. Actual water levels varied from approximately 620.5 feet amsl to 632.5 feet amsl. This represents an increase in the average water levels from the third quarter 2006. MWH will continue to monitor the water levels in both the SBPA and Off-Site Area to ensure vapor extraction at the ISVE wells is not inhibited.

7.0 SYSTEM OPERATION

The GWTP operated as designed for approximately 92 percent of the fourth quarter of 2006 (based on 2,014 hours of operation out of a total of 2,184 hours). The system drew influent water from the On-Site Area BWES, the Off-Site Area BWES, the PGCS, and MW-56.

The Off-Site Area ISVE system continued to operate as designed for approximately 74 percent of the fourth quarter of 2006 (based on 1,707 hours of operation out of a total of 2,184 hours). The SBPA ISVE system continued to operate as designed for approximately 78 percent of the fourth quarter of 2006 (based on 1,625 hours of operation out of a total of 2,184 hours).

A majority of the downtime for the ISVE systems was associated with maintenance of the thermal oxidizers.

8.0 CONCLUSIONS AND RECOMMENDATIONS

This section provides a summary of the operational status of the active remedial systems at the ACS NPL site for the subject period. Anticipated activities for the upcoming quarter and recommendations for system modifications are also provided.

8.1 GWTP OPERATION

The GWTP continued to operate normally during the fourth quarter of 2006. No significant modifications were made to the system during the period. The GWTP continued to treat water from all available sources, except MW-10C. The pump at MW-10C has malfunctioned and will be replaced.

The list of sources sending groundwater to the GWTP will be expanded upon the completion of the Lower Aquifer Pumping System. Due to the saturated ground conditions and the winter weather, MWH has suspended the installation of this system. MWH will monitor conditions in the area to determine an appropriate time to complete the installation.

Annual sediment sampling conducted in December 2007 indicated that concentrations of Aroclor-1248 were detected above laboratory method detection limits. Although concentrations remain below any action levels, results for this compound had been non-detect for previous sampling events. As a result, MWH will accelerate the schedule for collection of the 2007 sampling event to coincide with the next groundwater sampling event to be conducted in April 2007.

MWH is evaluating methods for using the free product regularly collected from the SBPA ISVE wells. Potential options include injection of the liquid into the combustion chamber of ThermOx 2 or using the material to fuel room heaters during the winter time.

8.2 ISVE OPERATON

The ISVE systems continued to operate normally during the fourth quarter of 2006. The operational times of both the systems were decreased primarily due to maintenance issues associated with the thermal oxidizers. MWH will continue to perform O&M services on these units to ensure adequate operational time for the ISVE systems. No significant changes were made to the operational configuration of the ISVE systems.

As shown in Figure 3.1, the VOC removal rates (in pounds per day) were observed to be within range of previous events. MWH is currently reviewing the equipment used to monitor the rate of VOC extraction by the ISVE systems. Alternate flow meters are being investigated to identify one that will allow MWH to more accurately and consistently measure the flow and mass of VOCs in the extracted vapor streams.

8.3 GROUNDWATER LEVEL MONITORING

As indicated in Section 6.0, the groundwater extraction system continues to successfully perform its intended function of isolating and protecting the groundwater outside the barrier wall from the source areas of the Site inside the barrier walls.

Recent groundwater level monitoring results indicate that levels in both the On-Site and Off-Site Areas have risen above previous minimum levels. MWH is evaluating the performance of the extraction trenches and wells to ensure correct operation and will continue to monitor groundwater levels.

8.4 HEALTH AND SAFETY

No health and safety incidents were reported during the fourth quarter of 2006. MWH continues to perform site activities in accordance with the site Health and Safety Plan and all applicable addendums.

MWH's annual safety meeting with ACS personnel was held on October 30, 2006. The meeting included an update on the status of MWH's work at the site and a refresher on the health and safety aspects of the ongoing operations. An update on ACS activities was provided for MWH personnel.

Eyewash stations were installed in both of the ISVE blower sheds. The annual fire extinguisher inspection was complete in November.

Health and Safety statistics for the ACS Site as of December 31, 2006 are:

- 3,508 consecutive days with no lost time due to an accident or Health and Safety incident.
- 1,200 consecutive days without an incident requiring first aid.

9.0 REFERENCES

1. *Final Remedial Design Report: Final Remedy, ACS NPL Site*, Montgomery Watson, August 1999.
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3. *Performance Standard Verification Plan, ACS NPL Site*, Montgomery Watson, June 1999.
4. *Phase I Technical Memorandum Wetland Investigation, ACS NPL Site*, Montgomery Watson, July 1996.
5. *Phase II Technical Memorandum Wetland Investigation, ACS NPL Site*, Montgomery Watson, February 1997.
6. *Quality Assurance Project Plan, ACS NPL Site*, Montgomery Watson Harza, March 2001.
7. *U.S. EPA Specifications and Guidance for Obtaining Contaminant-Free Sample Containers*, United States Environmental Protection Agency, 1992.
8. *Contract Laboratory Program National Functional Guidelines for Inorganic Data Review*, U.S. EPA, February 1994.
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TABLES

Table 2.1
Groundwater Treatment System Effluent Discharge Limits
American Chemical Service NPL Site
Griffith, Indiana

Groundwater Quality Parameter	Effluent Standard (Limit)
<i>General Water Quality Parameters</i>	
pH	6 - 9 S.U.
BOD-5	30 mg/L
TSS	30 mg/L
<i>Inorganics</i>	
Arsenic	50 µg/L
Beryllium	NE
Cadmium	4.1 µg/L
Manganese	NE
Mercury	0.02 µg/L (w/DL = 0.64)
Selenium	8.2 µg/L
Thallium	NE
Zinc	411 µg/L
<i>Volatile Organics</i>	
Acetone	6,800 µg/L
Benzene	5 µg/L
2-Butanone	210 µg/L
Chloromethane	NE
1,4 - Dichlorobenzene	NE
1,1 - Dichloroethane	NE
1,2 - Dichloroethene - cis	70 µg/L
Ethylbenzene	34 µg/L
Methylene chloride	5 µg/L
Tetrachloroethene	5 µg/L
Trichloroethene	5 µg/L
Vinyl chloride	2 µg/L
4 - Methyl - 2 - pentanone	15 µg/L
<i>Semi-Volatile Organics</i>	
bis(2 - Chloroethyl) ether	9.6 µg/L
bis(2 - Ethylhexyl) phthalate	6 µg/L
Isophorone	50 µg/L
4 - Methylphenol	34 µg/L
Pentachlorophenol	1 µg/L
<i>PCBs</i>	
PCBs	0.00056 µg/L (w/DL = 0.1 to 0.9)

Notes:

NE = No effluent limit established.

DL = Detection limit

S.U. = Standard pH units

µg/L - micrograms per Liter

Table 2.2
Summary of Effluent Analytical Results - Fourth Quarter 2006
Groundwater Treatment System
American Chemical Service NPL Site
Griffith, Indiana

Event Date	Month 113 10/11/2006	Month 114 11/1/2006	Month 115 12/20/2006	Effluent Limits	Lab Reporting Limits
pH	7.45 J	7.05 J	7.28 J	6-9	none
TSS	0.800 B/	NS	NS	30	10
BOD	< 2.0 B-01, U/	NS	NS	30	2
Arsenic	9.5 B/	NS	NS	50	3.4
Beryllium	0.45 B/UB	NS	NS	NE	0.2
Cadmium	0.20 U/	NS	NS	4.1	0.3
Manganese	0.10 U/	NS	NS	NE	10
Mercury	0.10 U/	NS	NS	0.02 (w/DL = 0.64)	0.64
Selenium	2.5 U/	NS	NS	8.2	4.3
Thallium	3.2 U/	NS	NS	NE	5.7
Zinc	1.1 B/UB	NS	NS	411	1.2
Benzene	0.50 U/UJ	0.50 U/	0.50 U/	5	0.5
Acetone	2.5 U/UJ	2.5 U/UJ	2.5 U/UJ	6,800	3
2-Butanone	2.5 U/UJ	2.5 U/	2.5 U/	210	3
Chloromethane	0.50 U/UJ	0.50 U/	0.50 U/UJ	NE	0.5
1,4-Dichlorobenzene	0.50 U/UJ	0.50 U/	0.50 U/	NE	0.5
1,1-Dichloroethane	0.50 U/UJ	0.15 J/	0.68 /	NE	0.5
cis-1,2-Dichloroethene	0.58 J/	0.42 J/	1.4 /	70	0.5
Ethylbenzene	0.50 U/UJ	0.50 U/	0.50 U/	34	0.5
Methylene chloride	0.45 J/J	0.84 /	0.65 /	5	0.6
Tetrachloroethene	0.50 U/UJ	0.50 U/	0.50 U/	5	0.5
Trichloroethene	0.50 U/UJ	0.50 U/	0.50 U/	5	0.5
Vinyl chloride	0.50 U/UJ	0.50 U/	0.40 J/J	2	0.5
4-Methyl-2-pentanone	2.5 U/UJ	2.5 U/	2.5 U/	15	3
bis (2-Chloroethyl) ether	ND	NS	NS	9.6	9.6
bis(2-Ethylhexyl) - phthalate	ND	NS	NS	6	6
4 - Methylphenol	ND	NS	NS	34	10
sophorone	ND	NS	NS	50	10
Pentachlorophenol	1.00 U/	NS	NS	1	1
PCB/Aroclor-1016	ND	NS	NS	0.00056 (w/DL = 0.1 to 0.9)	0.5
PCB/Aroclor-1221	ND	NS	NS	0.00056 (w/DL = 0.1 to 0.9)	0.92*
PCB/Aroclor-1232	ND	NS	NS	0.00056 (w/DL = 0.1 to 0.9)	0.5
PCB/Aroclor-1242	ND	NS	NS	0.00056 (w/DL = 0.1 to 0.9)	0.5
PCB/Aroclor-1248	ND	NS	NS	0.00056 (w/DL = 0.1 to 0.9)	0.5
PCB/Aroclor-1254	ND	NS	NS	0.00056 (w/DL = 0.1 to 0.9)	0.5
PCB/Aroclor-1260	ND	NS	NS	0.00056 (w/DL = 0.1 to 0.9)	0.5

Notes:

J= folded result indicates a exceedence of the discharge limit

pH data is expressed in S.U.

Metals, VOC, SVOC and PCB data is expressed in ug/L

ND = Not detected

NS = This analyte was not sampled or analyzed for

NE = No effluent limit established.

DL = Detection limit

.. = Approved SW-846 method is incapable of achieving effluent limit.

Suffix Definitions:

-' = Data qualifier added by laboratory

/_ = Data qualifier added by data validator

J = Result is detected below the reporting limit and is an estimated concentration

U = Analyte is not detected at or above the indicated concentration

B = Compound is also detected in the blank

UJ = Indicates the compound or analyte was analyzed for but not detected. The sample detection limit is an estimated value

UB = Compound or analyte is not detected at or above the indicated concentration due to blank contamination

B-01 = Sample dilutions set-up for BOD analysis did not meet the oxygen depletion criteria of at least 2mg/L dissolved oxygen depletion.

Therefore, the reported result is an estimated value only.

Table 2.3
Summary of Sediment Analytical Results
Groundwater Treatment System
American Chemical Service NPL Site
Griffith, Indiana

PCB Compound	Results (ug/kg)														
	12/4/1998	2/3/2000	2/3/00 DUP	8/21/2001	8/21/01 DUP	6/5/2002	6/5/02 DUP	1/13/04	1/13/04 DUP	9/27/04	9/27/04 DUP	6/15/05	6/15/05 DUP	12/11/06	12/11/06 DUP
Aroclor-1016	ND (33)	ND (59)	ND (79)	ND (62) /UJ	ND (71)	ND (52) /UJ	ND (49)	ND (67)	ND (76)	ND (62)	ND (78)	ND (74)	ND (39)	ND (71)	ND (53)
Aroclor-1221	ND (33)	ND (77)	ND (100)	ND (82) /UJ	ND (92)	ND (67) /UJ	ND (64)	ND (84)	ND (95)	ND (84)	ND (110)	ND (100)	ND (53)	ND (100)	ND (76)
Aroclor-1232	ND (33)	ND (59)	ND (79)	ND (62) /UJ	ND (71)	ND (52) /UJ	ND (49)	ND (67)	ND (76)	ND (62)	ND (78)	ND (74)	ND (39)	ND (50)	ND (38)
Aroclor-1242	ND (33)	ND (41)	ND (55)	ND (43) /UJ	ND (49) /UJ	ND (36) /UJ	ND (34)	ND (42)	ND (48)	ND (42)	ND (53)	ND (50)	ND (27)	ND (50)	ND (38)
Aroclor-1248	ND (33)	ND (41)	ND (55)	ND (43) /UJ	ND (49) /UJ	ND (36) /UJ	ND (34)	ND (42)	ND (48)	ND (42)	ND (53)	ND (49)	ND (27)	300	450
Aroclor-1254	ND (33)	22 J/	15 J/	73 P/J	39 JP/J	ND (36) /UJ	ND (34)	ND (42)	ND (48)	ND (42)	ND (53)	ND (54)	ND (27)	ND (50)	ND (38)
Aroclor-1260	ND (33)	ND (59)	ND (79)	ND (62) /UJ	ND (71) /UJ	41 J/J	ND (49)	35 J	ND (76)	ND (62)	ND (78)	ND (74)	ND (39)	ND (50)	ND (38)
Total PCBs⁴	ND	22	15	73	39	41	ND	35	ND	ND	ND	ND	ND	300	450

Notes:

1 ND () = Compound was not detected. The detection limit is included in parentheses

2 December 4, 1998 sample was analyzed by Quanterra. All other samples were analyzed by Compuchem.

3 DUP = Duplicate sample

4 The total PCB value presented here are estimated totals based on estimated concentrations of individual Aroclors

Suffix Definitions:

/ = Data qualifier added by laboratory

/_ = Data qualifier added by data validator

J = Result is detected below the reporting limit and is an estimated concentration

P = The Relative Percent Difference (RPD) between the two GC column values is greater than 25%. The higher value has been reported.

JP = The Relative Percent Difference (RPD) between the two GC column values is greater than 25%. The higher value has been reported. The concentration is also estimated.

UJ = Indicates the compound or analyte was analyzed for but not detected. The sample detection limit is an estimated value.

Table 3.1
Thermal Oxidizer 1 Results for Method TO-14 (VOCs) - October 2006
American Chemical Service
Griffith, Indiana

Compounds	Units	Sampled 10/19/06							
		Therm-Ox 1				Destruction Efficiency			
		Influent	Influent Dup	Effluent	Low	High	Average		
1,1,-Trichloroethane	ppbv	10,000	/R	10,000	/R	35	/R	99.65%	99.65% 99.65%
1,1,2,2-Tetrachloroethane	ppbv	94	J/R	ND	U/R	0.76	/R	NC	NC NC
1,1,2-Trichloroethane	ppbv	ND	U/R	ND	U/R	ND	U/R	NC	NC NC
1,1-Dichloroethane	ppbv	1,400	/R	1,400	/R	5.3	/R	99.62%	99.62% 99.62%
1,1-Dichloroethene	ppbv	240	/R	220	/R	57	/R	74.09%	76.25% 75.17%
1,2-Dichloroethane	ppbv	210	/R	220	/R	1.1	/R	99.48%	99.50% 99.49%
1,2-Dichloropropane	ppbv	230	/R	190	/R	0.76	/R	99.60%	99.67% 99.63%
2-Butanone (Methyl Ethyl Ketone)	ppbv	410	J/R	390	J/R	14	/R	NC	NC NC
2-Hexanone	ppbv	ND	U/R	ND	U/R	0.69	J/R	NC	NC NC
4-Methyl-2-pentanone	ppbv	1,400	/R	1,600	/R	17	/R	98.79%	98.94% 98.86%
Acetone	ppbv	800	/R	620	/R	85	/R	86.29%	89.38% 87.83%
Benzene	ppbv	4,300	/R	4,000	/R	72	/R	98.20%	98.33% 98.26%
Bromodichloromethane	ppbv	ND	U/R	ND	U/R	ND	U/R	NC	NC NC
Bromoform	ppbv	ND	U/R	ND	U/R	ND	U/R	NC	NC NC
Bromomethane	ppbv	ND	U/R	ND	U/R	ND	U/R	NC	NC NC
Carbon Disulfide	ppbv	ND	U/R	ND	U/R	1.2	J/R	NC	NC NC
Carbon Tetrachloride	ppbv	ND	U/R	ND	U/R	0.55	J/R	NC	NC NC
Chlorobenzene	ppbv	ND	U/R	58	J/R	1.9	/R	NC	NC NC
Chloroethane	ppbv	170	/R	220	/R	1.9	/R	98.88%	99.14% 99.01%
Chloroform	ppbv	3,500	/R	3,400	/R	16	/R	99.53%	99.54% 99.54%
Chloromethane	ppbv	ND	U/R	ND	U/R	9	/R	NC	NC NC
cis-1,2-Dichloroethene	ppbv	9,000	/R	9,700	/R	58	/R	99.36%	99.40% 99.38%
cis-1,3-Dichloropropene	ppbv	ND	U/R	ND	U/R	0.21	J/R	NC	NC NC
Dibromochloromethane	ppbv	ND	U/R	ND	U/R	ND	U/R	NC	NC NC
Ethyl Benzene	ppbv	8,300	/R	8,000	/R	32	/R	99.60%	99.61% 99.61%
m,p-Xylene	ppbv	33,000	/R	32,000	/R	130	/R	99.59%	99.61% 99.60%
Methylene Chloride	ppbv	4,700	/R	5,000	/R	170	/R	96.38%	96.60% 96.49%
o-Xylene	ppbv	13,000	/R	13,000	/R	54	/R	99.58%	99.58% 99.58%
Styrene	ppbv	ND	U/R	ND	U/R	17	/R	NC	NC NC
Tetrachloroethene	ppbv	26,000	/R	26,000	/R	220	/R	99.15%	99.15% 99.15%
Toluene	ppbv	34,000	/R	34,000	/R	130	/R	99.62%	99.62% 99.62%
trans-1,2-Dichloroethene	ppbv	170	J/R	110	J/R	14	/R	NC	NC NC
trans-1,3-Dichloropropene	ppbv	ND	U/R	ND	U/R	ND	U/R	NC	NC NC
Trichloroethene	ppbv	9,100	/R	9,300	/R	72	/R	99.21%	99.23% 99.22%
Vinyl Chloride	ppbv	1,500	/R	1,800	/R	30	/R	98.00%	98.33% 98.17%
Total	ppbv	161,524		161,228		1,246.37		99.23%	99.23% 99.23%
Total	lb/hr	2.118		2.112		0.015		99.29%	99.29% 99.29%

Notes:

NC - Not calculated

ND - Non-detected

ppbv - parts per billion volume

lb/hr - pounds per hour

Qualifiers:

J - Result is estimated

U - below reported quantitation limit

R - Quality control indicates the data is not usable

/ - Laboratory data qualifier

/ - Data validation qualifier

Destruction efficiencies were not calculated if either the influent or effluent samples were estimated.

Destruction efficiencies were also not calculated if the effluent result exceeded either influent result.

Total destruction efficiencies that include the estimated results of any individual compound will be considered an estimated value.

System	Date	Influent Temp (°F)	Effluent Temp (°F)	Flow (scfm)
SBPA	10/19/06	96	142	1420

Table 3.2
Thermal Oxidizer 1 Results for Method TO-14 (VOCs) - November 2006
American Chemical Service
Griffith, Indiana

Compounds	Units	Sampled 11/2/06								
		Therm-Ox 1				Destruction Efficiency				
		Influent	Influent Dup	Effluent		Low	High	Average		
1,1,1-Trichloroethane	ppbv	18,000		16,000	3.4		99.98%	99.98%	99.98%	
1,1,2-Tetrachloroethane	ppbv	ND	U	ND	U	ND	NC	NC	NC	
1,1,2-Trichloroethane	ppbv	51	J/J	56	J/J	ND	U	NC	NC	
1,1-Dichloroethane	ppbv	2,100		2,200		1.5		99.93%	99.93%	99.93%
1,1-Dichloroethene	ppbv	1,100		840		210		75.00%	80.91%	77.95%
1,2-Dichloroethane	ppbv	300		260		0.81	J/J	NC	NC	NC
1,2-Dichloropropane	ppbv	250		250		ND	U	100.00%	100.00%	100.00%
2-Butanone (Methyl Ethyl Ketone)	ppbv	580		530	J/J	8	NC	NC	NC	
2-Fexanone	ppbv	ND	U	ND	U	ND	U	NC	NC	NC
4-Methyl-2-pentanone	ppbv	660		880		5.2		99.21%	99.41%	99.31%
Acetone	ppbv	830		910		24		97.11%	97.36%	97.24%
Benzene	ppbv	2,700		3,200		170		93.70%	94.69%	94.20%
Bromodichloromethane	ppbv	ND	U	ND	U	0.36	J/J	NC	NC	NC
Bromoform	ppbv	ND	U	ND	U	ND	U	NC	NC	NC
Bromomethane	ppbv	ND	U	ND	U	ND	U	NC	NC	NC
Carbon Disulfide	ppbv	ND	U	ND	U	5.4		NC	NC	NC
Carbon Tetrachloride	ppbv	ND	U	ND	U	1.1		NC	NC	NC
Chlorobenzene	ppbv	ND	U	ND	U	8.8		NC	NC	NC
Chloroethane	ppbv	180		500		1.6		99.11%	99.68%	99.40%
Chloroform	ppbv	4,600		4,300		7.4		99.83%	99.84%	99.83%
Chloromethane	ppbv	ND	U	ND	U	15		NC	NC	NC
cis-1,2-Dichloroethene	ppbv	11,000		13,000		98		99.11%	99.25%	99.18%
cis-1,3-Dichloropropene	ppbv	ND	U	ND	U	0.81	J/J	NC	NC	NC
Dibromochloromethane	ppbv	ND	U	ND	U	ND	U	NC	NC	NC
Ethyl Benzene	ppbv	5,600		6,000		6		99.89%	99.90%	99.90%
m,p-Xylene	ppbv	23,000		24,000		30		99.87%	99.88%	99.87%
Methylene Chloride	ppbv	6,500		5,600		50		99.11%	99.23%	99.17%
o-Xylene	ppbv	11,000		10,000		11		99.89%	99.90%	99.90%
Styrene	ppbv	ND	U	ND	U	25		NC	NC	NC
Tetrachloroethene	ppbv	26,000		24,000		390		98.38%	98.50%	98.44%
Toluene	ppbv	31,000		33,000		35		99.89%	99.89%	99.89%
trans-1,2-Dichloroethene	ppbv	95	J/J	100	J/J	61		NC	NC	NC
trans-1,3-Dichloropropene	ppbv	ND	U	ND	U	0.72	J/J	NC	NC	NC
Trichloroethene	ppbv	15,000		13,000		160		98.77%	98.93%	98.85%
Vinyl Chloride	ppbv	1,200		1,800		91		92.42%	94.94%	93.68%
Total	ppbv	161,746		160,426		1,421.10		99.11%	99.12%	99.12%
Total	lb/hr	4.316		4.212		0.037		99.12%	99.14%	99.13%

Notes:

NC - Not calculated

ND - Non-detect

ppbv - parts per billion volume

lb/hr - pounds per hour

Destruction efficiencies were not calculated if either the influent or effluent samples were estimated.

Destruction efficiencies were also not calculated if the effluent result exceeded either influent result.

Total destruction efficiencies that include the estimated results of any individual compound will be considered an estimated value.

Qualifiers:

J - Result is estimated

U - below reported quantitation limit

/ - Laboratory data qualifier

/ - Data validation qualifier

System	Date	Influent Temp (°F)	Effluent Temp (°F)	Flow (scfm)
SBPA	11/02/06	98	130	1693

Table 3.3
Thermal Oxidizer 1 Results for Method TO-14 (VOCs) - December 2006
American Chemical Service
Griffith, Indiana

Compounds	Units	Sampled 12/11/06							
		Therm-Ox 1				Destruction Efficiency			
		Influent	Influent Dup	Effluent		Low	High	Average	
1,1,1-Trichloroethane	ppbv	18,000		43		99.76%	99.77%	99.77%	
1,1,2,2-Tetrachloroethane	ppbv	ND	U	ND	U	NC	NC	NC	
1,1,2-Trichloroethane	ppbv	47	J/J	70	J/J	0.25	J/J	NC	NC
1,1-Dichloroethane	ppbv	3,100		7.9		99.75%	99.76%	99.75%	
1,1-Dichloroethene	ppbv	1,200		200		83.33%	84.62%	83.97%	
1,2-Dichloroethane	ppbv	270		0.85		99.69%	99.70%	99.69%	
1,2-Dichloropropane	ppbv	300		0.77		99.74%	99.77%	99.76%	
2-Butanone (Methyl Ethyl Ketone)	ppbv	940		560	E	NC	NC	NC	
2-Hexanone	ppbv	ND	U	ND	U	1.2	J/J	NC	NC
4-Methyl-2-pentanone	ppbv	1,500		44		97.07%	97.56%	97.31%	
Acetone	ppbv	1,200		23		98.08%	98.23%	98.16%	
Benzene	ppbv	4,800		120		97.50%	97.86%	97.68%	
Bromodichloromethane	ppbv	ND	U	ND	U	NC	NC	NC	
Bromoform	ppbv	ND	U	ND	U	NC	NC	NC	
Bromomethane	ppbv	ND	U	ND	U	NC	NC	NC	
Carbon Disulfide	ppbv	33	J/J	39	J/J	120		NC	NC
Carbon Tetrachloride	ppbv	ND	U	ND	U	0.6	J/J	NC	NC
Chlorobenzene	ppbv	69	J/J	97	J/J	6.3		NC	NC
Chloroethane	ppbv	770	J/J	780	J/J	3.2	J/J	99.58%	99.59%
Chloroform	ppbv	4,900		5,300		13		99.73%	99.75%
Chloromethane	ppbv	ND	U	ND	U	8.9		NC	NC
cis 1,2-Dichloroethene	ppbv	15,000		17,000		87		99.42%	99.49%
cis 1,3-Dichloropropene	ppbv	ND	U	ND	U	0.25	J/J	NC	NC
Dibromochloromethane	ppbv	ND	U	ND	U	NC	NC	NC	
Ethyl Benzene	ppbv	9,300		11,000		24		99.74%	99.78%
m,p-Xylene	ppbv	39,000		46,000		110		99.72%	99.76%
Methylene Chloride	ppbv	6,600		7,500		50		99.24%	99.33%
o-Xylene	ppbv	18,000		21,000		50		99.72%	99.76%
Styrene	ppbv	ND	U	ND	U	12		NC	NC
Tetrachloroethene	ppbv	26,000		28,000		260		99.00%	99.07%
Toluene	ppbv	41,000		48,000		100		99.76%	99.79%
trans-1,2-Dichloroethene	ppbv	110	J/J	140	J/J	43		NC	NC
trans-1,3-Dichloropropene	ppbv	ND	U	ND	U	0.24	J/J	NC	NC
Trichloroethene	ppbv	14,000		15,000		120		99.14%	99.20%
Vinyl Chloride	ppbv	1,500		1,500		76		94.93%	94.93%
Total	ppbv	207,639		235,446		2,085.46		99.00%	99.11%
Total	lb/hr	5.359		6.040		0.047		99.12%	99.22%
									99.17%

Notes:

NC - Not calculated

ND - Non-detect

ppbv - parts per billion volume

lb/hr - pounds per hour

Qualifiers:

J - Result is estimated

U - below reported quantitation limit

E - Exceeds instrument calibration range

/ - Laboratory data qualifier

/_ - Data validation qualifier

Destruction efficiencies were not calculated if either the influent or effluent samples were estimated.

Destruction efficiencies were also not calculated if the effluent result exceeded either influent result.

Total destruction efficiencies that include the estimated results of any individual compound will be considered an estimated value.

System	Date	Influent Temp (°F)	Effluent Temp (°F)	Flow (scfm)
SBPA	12/11/06	100	142	1707

Table 3.4
Thermal Oxidizer 2 Results for Method TO-14 (VOCs) - October 2006
American Chemical Service
Griffith, Indiana

Compounds	Units	Sampled 10/19/06								
		Therm-Ox 2				Destruction Efficiency				
		Influent	Influent Dup	Effluent		Low	High	Average		
1,1,1-Trichloroethane	ppbv	29,000	/R	26,000	/R	430	/R	98.35%	98.52%	98.43%
1,1,2,2-Tetrachloroethane	ppbv	ND	U/R	ND	U/R	ND	U/R	NC	NC	NC
1,1,2-Trichloroethane	ppbv	ND	U/R	ND	U/R	ND	U/R	NC	NC	NC
1,1-Dichloroethane	ppbv	3,800	/R	3,600	/R	60	/R	98.33%	98.42%	98.38%
1,1-Dichloroethene	ppbv	300	J/R	350	/R	160	/R	NC	NC	NC
1,2-Dichloroethane	ppbv	820	/R	780	/R	13	/R	98.33%	98.41%	98.37%
1,2-Dichloropropane	ppbv	290	J/R	240	J/R	3.8	J/R	NC	NC	NC
2-Butanone (Methyl Ethyl Ketone)	ppbv	12,000	/R	12,000	/R	200	/R	98.33%	98.33%	98.33%
2-Hexanone	ppbv	ND	U/R	ND	U/R	2.6	J	NC	NC	NC
4-Methyl-2-pentanone	ppbv	6,600	/R	6,100	/R	61	/R	99.00%	99.08%	99.04%
Acetone	ppbv	16,000	/R	19,000	/R	470	/R	97.06%	97.53%	97.29%
Benzene	ppbv	15,000	/R	15,000	/R	480	/R	96.80%	96.80%	96.80%
Bromodichloromethane	ppbv	ND	U/R	ND	U/R	ND	U/R	NC	NC	NC
Bromoform	ppbv	ND	U/R	ND	U/R	ND	U/R	NC	NC	NC
Bromomethane	ppbv	ND	U/R	ND	U/R	ND	U/R	NC	NC	NC
Carbon Disulfide	ppbv	210	J/R	ND	U/R	9.2	J/R	NC	NC	NC
Carbon Tetrachloride	ppbv	ND	U/R	ND	U/R	3.8	J/R	NC	NC	NC
Chlorobenzene	ppbv	ND	U/R	ND	U/R	4	J/R	NC	NC	NC
Chloroethane	ppbv	570	/R	540	/R	10	/R	98.15%	98.25%	98.20%
Chloroform	ppbv	1,800	/R	1,900	/R	36	/R	98.00%	98.11%	98.05%
Chloromethane	ppbv	230	J/R	ND	U/R	ND	U/R	NC	NC	NC
cis-1,2-Dichloroethene	ppbv	6,000	/R	6,800	/R	120	/R	98.00%	98.24%	98.12%
cis-1,3-Dichloropropene	ppbv	ND	U/R	ND	U/R	ND	U/R	NC	NC	NC
Dibromochloromethane	ppbv	ND	U/R	ND	U/R	ND	U/R	NC	NC	NC
Ethyl Benzene	ppbv	9,800	/R	12,000	/R	120	/R	98.78%	99.00%	98.89%
m,p-Xylene	ppbv	37,000	/R	51,000	/R	430	/R	98.84%	99.16%	99.00%
Methylene Chloride	ppbv	30,000	/R	30,000	/R	580	/R	98.07%	98.07%	98.07%
o-Xylene	ppbv	13,000	/R	18,000	/R	150	/R	98.85%	99.17%	99.01%
Styrene	ppbv	ND	U/R	ND	U/R	95	/R	NC	NC	NC
Tetrachloroethene	ppbv	23,000	/R	24,000	/R	570	/R	97.52%	97.63%	97.57%
Toluene	ppbv	89,000	/R	100,000	/R	1,400	/R	98.43%	98.60%	98.51%
trans-1,2-Dichloroethene	ppbv	ND	U/R	ND	U/R	22.0	/R	NC	NC	NC
trans-1,3-Dichloropropene	ppbv	ND	U/R	ND	U/R	ND	U/R	NC	NC	NC
Trichloroethene	ppbv	13,000	/R	14,000	/R	290	/R	97.77%	97.93%	97.85%
Vinyl Chloride	ppbv	1,500	/R	1,400	/R	57	/R	95.93%	96.20%	96.06%
Total	ppbv	308,920		342,710		5,777.4		98.13%	98.31%	98.22%
Total	lb/hr	7.713		8.522		0.1429		98.15%	98.32%	98.24%

Notes:

NC - Not calculated

ND - Non-detect

ppbv - parts per billion volume

lb/hr - pounds per hour

Qualifiers:

J - Result is estimated

U - below reported quantitation limit

R - Quality control indicates the data is not usable

/ - Laboratory data qualifier

/_ - Data validation qualifier

Destruction efficiencies were not calculated if either the influent or effluent samples were estimated.

Destruction efficiencies were also not calculated if the effluent result exceeded either influent result.

Total destruction efficiencies that include the estimated results of any individual compound will be considered an estimated value.

System	Date	Influent Temp (°F)	Effluent Temp (°F)	Flow (scfm)
Off-Site	10/19/06	72	NA	1707

Table 3.5
Thermal Oxidizer 2 Results for Method TO-14 (VOCs) - November 2006
American Chemical Service
Griffith, Indiana

Compounds	Units	Sampled 11/2/06						
		Therm-Ox 2			Destruction Efficiency			
		Influent	Influent Dup	Effluent	Low	High	Average	
1,1,1-Trichloroethane	ppbv	20,000		460	97.70%	97.70%	97.70%	
1,1,2,2-Tetrachloroethane	ppbv	ND	U	ND	NC	NC	NC	
1,1,2-Trichloroethane	ppbv	130	J/J	140	J/J	NC	NC	NC
1,1-Dichloroethane	ppbv	2,800		62	97.79%	97.79%	97.79%	
1,1-Dichloroethene	ppbv	820		180	78.05%	80.22%	79.13%	
1,2-Dichloroethane	ppbv	650		16	97.46%	97.54%	97.50%	
1,2-Dichloropropane	ppbv	160	J/J	210	J/J	NC	NC	NC
2-Butanone (Methyl Ethyl Ketone)	ppbv	16,000		990	92.38%	93.81%	93.10%	
2-Hexanone	ppbv	ND	U	8.2	J/J	NC	NC	NC
4-Methyl-2-pentanone	ppbv	6,000		210	96.38%	96.50%	96.44%	
Acetone	ppbv	15,000		1100	89.00%	92.67%	90.83%	
Benzene	ppbv	11,000		490	95.55%	95.92%	95.73%	
Bromodichloromethane	ppbv	ND	U	ND	NC	NC	NC	
Bromoform	ppbv	ND	U	ND	NC	NC	NC	
Bromomethane	ppbv	ND	U	ND	NC	NC	NC	
Carbon Disulfide	ppbv	ND	U	ND	NC	NC	NC	
Carbon Tetrachloride	ppbv	ND	U	4.1	J/J	NC	NC	NC
Chlorobenzene	ppbv	ND	U	ND	4.3	J/J	NC	NC
Chloroethane	ppbv	ND	U	ND	5.6	J/J	NC	NC
Chloroform	ppbv	1,700		46	97.29%	97.29%	97.29%	
Chloromethane	ppbv	ND	U	17	J/J	NC	NC	NC
cis-1,2-Dichloroethene	ppbv	4,900		140	97.02%	97.14%	97.08%	
cis-1,3-Dichloropropene	ppbv	ND	U	ND	NC	NC	NC	
Dibromochloromethane	ppbv	ND	U	ND	NC	NC	NC	
Ethyl Benzene	ppbv	9,200		180	98.04%	98.20%	98.12%	
m,p-Xylene	ppbv	38,000		700	98.16%	98.41%	98.28%	
Methylene Chloride	ppbv	15,000		430	97.13%	97.13%	97.13%	
o-Xylene	ppbv	14,000		270	98.07%	98.31%	98.19%	
Styrene	ppbv	ND	U	ND	57	NC	NC	NC
Tetrachloroethene	ppbv	16,000		640	96.00%	96.24%	96.12%	
Toluene	ppbv	85,000		1800	97.88%	98.04%	97.96%	
trans-1,2-Dichloroethene	ppbv	ND	U	ND	NC	NC	NC	
trans-1,3-Dichloropropene	ppbv	ND	U	ND	NC	NC	NC	
Trichloroethene	ppbv	13,000		380	97.08%	97.29%	97.18%	
Vinyl Chloride	ppbv	550		54	89.62%	90.18%	89.90%	
Total	ppbv	269,910		8,267.8	96.94%	97.05%	96.99%	
Total	lb/hr	4.010		0.1170	97.08%	97.23%	97.15%	

Notes:

NC - Not calculated

ND - Non-detect

ppbv - parts per billion volume

lb/hr - pounds per hour

Destruction efficiencies were not calculated if either the influent or effluent samples were estimated.

Destruction efficiencies were also not calculated if the effluent result exceeded either influent result.

Total destruction efficiencies that include the estimated results of any individual compound will be considered an estimated value.

Qualifiers:

J - Result is estimated

U - below reported quantitation limit

/ - Laboratory data qualifier

_ - Data validation qualifier

System	Date	Influent Temp (°F)	Effluent Temp (°F)	Flow (scfm)
Off-Site	11/02/06	68	NA	1370

Table 3.6
Thermal Oxidizer 2 Results for Method TO-14 (VOCs) - December 2006
American Chemical Service
Griffith, Indiana

Compounds	Units	Sampled 12/11/06						
		Therm-Ox 2			Destruction Efficiency			
		Influent	Influent Dup	Effluent	Low	High	Average	
,1,1-Trichloroethane	ppbv	22,000		720	96.57%	96.73%	96.65%	
,1,2,2-Tetrachloroethane	ppbv	ND	U	2.8	J/J	NC	NC	NC
,1,2-Trichloroethane	ppbv	140	J/J	5.4	J/J	NC	NC	NC
,1-Dichloroethane	ppbv	2,800		110	95.93%	96.07%	96.00%	
,1-Dichloroethene	ppbv	380		190	50.00%	55.81%	52.91%	
,2-Dichloroethane	ppbv	650		25	96.15%	96.15%	96.15%	
,2-Dichloropropane	ppbv	200	J/J	7.2	J/J	NC	NC	NC
-Butanone (Methyl Ethyl Ketone)	ppbv	16,000		500	75.00%	96.88%	85.94%	
-Hexanone	ppbv	180	J/J	5.7	J/J	NC	NC	NC
4-Methyl-2-pentanone	ppbv	6,600		120	95.56%	98.18%	96.87%	
Acetone	ppbv	15,000		770	90.83%	94.87%	92.85%	
Benzene	ppbv	13,000		820	93.17%	93.69%	93.43%	
Bromodichloromethane	ppbv	ND	U	ND	U	NC	NC	NC
Bromoform	ppbv	ND	U	ND	U	NC	NC	NC
Bromomethane	ppbv	ND	U	ND	U	NC	NC	NC
Carbon Disulfide	ppbv	170	J/J	140	J/J	NC	NC	NC
Carbon Tetrachloride	ppbv	ND	U	ND	U	NC	NC	NC
Chlorobenzene	ppbv	ND	U	ND	U	NC	NC	NC
Chloroethane	ppbv	ND	U	ND	U	NC	NC	NC
Chloroform	ppbv	1,800		77	95.72%	95.72%	95.72%	
Chloromethane	ppbv	ND	U	18	J/J	NC	NC	NC
cis-1,2-Dichloroethene	ppbv	1,400		81	93.77%	94.21%	93.99%	
cis-1,3-Dichloropropene	ppbv	ND	U/R	ND	U/R	NC	NC	NC
Dibromochloromethane	ppbv	ND	U	ND	U	NC	NC	NC
Ethyl Benzene	ppbv	9,700		270	97.03%	97.22%	97.12%	
m,p-Xylene	ppbv	40,000		1100	97.11%	97.25%	97.18%	
Methylene Chloride	ppbv	19,000		870	95.17%	95.42%	95.29%	
c-Xylene	ppbv	15,000		400	97.14%	97.33%	97.24%	
Styrene	ppbv	ND	U	150	NC	NC	NC	
Tetrachloroethene	ppbv	16,000		780	95.13%	95.13%	95.13%	
Toluene	ppbv	83,000		2900	96.33%	96.51%	96.42%	
trans-1,2-Dichloroethene	ppbv	ND	U	12	J/J	NC	NC	NC
trans-1,3-Dichloropropene	ppbv	ND	U	ND	U	NC	NC	NC
Trichloroethene	ppbv	13,000		550	95.42%	95.77%	95.59%	
Vinyl Chloride	ppbv	ND	U	40	NC	NC	NC	
Total	ppbv	276,020	239,530	10,537.4	95.60%	96.18%	95.89%	
Total	lb/hr	4.104	3.663	0.155	95.77%	96.22%	96.00%	

Notes:

NC - Not calculated

ND - Non-detect

ppbv - parts per billion volume

lb/hr - pounds per hour

Qualifiers:

J - Result is estimated

U - below reported quantitation limit

R - Quality control indicates the data is not usable

/ - Laboratory data qualifier

/_ - Data validation qualifier

Destruction efficiencies were not calculated if either the influent or effluent samples were estimated.

Destruction efficiencies were also not calculated if the effluent result exceeded either influent result.

Total destruction efficiencies that include the estimated results of any individual compound will be considered an estimated value.

System	Date	Influent Temp (°F)	Effluent Temp (°F)	Flow (scfm)
Off-Site	39062	66	NA	1090

Table 3.7
SBPA and Off-Site ISVE System Results
for Method TO-14 (VOCs) - October 2006
American Chemical Service
Griffith, Indiana

Compounds	Units	Sampled 10/19/2006			
		SBPA ISVE		Off-Site ISVE	
1,1,1-Trichloroethane	ppbv	9,700	/R	36,000	/R
1,1,2,2-Tetrachloroethane	ppbv	ND	U/R	ND	U/R
1,1,2-Trichloroethane	ppbv	ND	U/R	ND	U/R
1,1-Dichloroethane	ppbv	1,400	/R	4,600	/R
1,1-Dichloroethene	ppbv	230	/R	350	J/R
1,2-Dichloroethane	ppbv	240	/R	900	/R
1,2-Dichloropropane	ppbv	210	/R	320	J/R
2-Butanone (Methyl Ethyl Ketone)	ppbv	420	J/J	16,000	/R
2-Hexanone	ppbv	ND	U/R	280.0	J/R
4-Methyl-2-pantanone	ppbv	1,700	/R	9,500	/R
Acetone	ppbv	750	/R	24,000	/R
Benzene	ppbv	4,700	/R	18,000	/R
Bromodichloromethane	ppbv	ND	U/R	ND	U/R
Bromoform	ppbv	ND	U/R	ND	U/R
Bromomethane	ppbv	ND	U/R	ND	U/R
Carbon Disulfide	ppbv	ND	U/R	180	J/R
Carbon Tetrachloride	ppbv	ND	U/R	ND	U/R
Chlorobenzene	ppbv	59	J/J	ND	U/R
Chloroethane	ppbv	160	/R	ND	U/R
Chloroform	ppbv	3,600	/R	2,500	/R
Chloromethane	ppbv	ND	U/R	ND	U/R
cis-1,2-Dichloroethene	ppbv	8,700	/R	1,400	/R
cis-1,3-Dichloropropene	ppbv	ND	U/R	ND	U/R
Dibromochloromethane	ppbv	ND	U/R	ND	U/R
Ethyl Benzene	ppbv	8,100	/R	13,000	/R
m,p-Xylene	ppbv	33,000	/R	50,000	/R
Methylene Chloride	ppbv	4,800	/R	44,000	/R
o-Xylene	ppbv	12,000	/R	18,000	/R
Styrene	ppbv	ND	U/R	ND	U/R
Tetrachloroethene	ppbv	24,000	/R	26,000	/R
Toluene	ppbv	35,000	/R	120,000	/R
trans-1,2-Dichloroethene	ppbv	ND	U/R	ND	U/R
trans-1,3-Dichloropropene	ppbv	ND	U/R	ND	U/R
Trichloroethene	ppbv	9,300	/R	16,000	/R
Vinyl Chloride	ppbv	1,400	/R	330	J/R
Total	ppbv	159,469		401,360	
Total	lb/hr	2.076		9.919	

Notes:

ND - Non-detect

ppbv - parts per billion volume

lb/hr - pounds per hour

Qualifiers:

J - Result is estimated

U - below reported quantitation limit

R - Quality control indicates the data is not usable

/ - Laboratory data qualifier

/ - Data validation qualifier

System	Date	Temp (°F)	Flow (scfm)
Off-Site	10/19/06	25	1707
SBPA	10/19/06	96	1420

Table 3.8
SBPA and Off-Site ISVE System Results
for Method TO-14 (VOCs) - November 2006
American Chemical Service
Griffith, Indiana

Compounds	Units	Sampled 11/2/2006			
		SBPA ISVE		Off-Site ISVE	
1,1,1-Trichloroethane	ppbv	18,000		25,000	
1,1,2,2-Tetrachloroethane	ppbv	ND	U	ND	U
1,1,2-Trichloroethane	ppbv	ND	U	180	J/J
1,1-Dichloroethane	ppbv	2,300		3,100	
1,1-Dichloroethene	ppbv	1,100		1,900	
1,2-Dichloroethane	ppbv	290		840	
1,2-Dichloropropane	ppbv	270		220	J/J
2-Butanone (Methyl Ethyl Ketone)	ppbv	810		16,000	
2-Hexanone	ppbv	ND	U	ND	U
4-Methyl-2-pentanone	ppbv	1,000		7,500	
Acetone	ppbv	1,300		13,000	
Benzene	ppbv	3,500		13,000	
Bromodichloromethane	ppbv	ND	U	ND	U
Eromoform	ppbv	ND	U	ND	U
Eromomethane	ppbv	ND	U	ND	U
Carbon Disulfide	ppbv	ND	U	ND	U
Carbon Tetrachloride	ppbv	ND	U	ND	U
Chlorobenzene	ppbv	ND	U	ND	U
Chloroethane	ppbv	380		ND	U
Chloroform	ppbv	4,700		2,200	
Chloromethane	ppbv	ND	U	ND	U
cis-1,2-Dichloroethene	ppbv	14,000		1,600	
cis-1,3-Dichloropropene	ppbv	ND	U	ND	U
Dibromochloromethane	ppbv	ND	U	ND	U
Ethyl Benzene	ppbv	7,300		14,000	
m,p-Xylene	ppbv	29,000		57,000	
Methylene Chloride	ppbv	5,600		22,000	
c-Xylene	ppbv	13,000		21,000	
Styrene	ppbv	ND	U	ND	U
Tetrachloroethene	ppbv	26,000		22,000	
Toluene	ppbv	39,000		110,000	
trans-1,2-Dichloroethene	ppbv	110	J/J	ND	U
trans-1,3-Dichloropropene	ppbv	ND	U	ND	U
Trichloroethene	ppbv	14,000		17,000	
Vinyl Chloride	ppbv	1,800		ND	U
Total	ppbv	183,460		347,540	
Total	lb/hr	4.792		5.247	

Notes:

ND - Non-detect

ppbv - parts per billion volume

lb/hr - pounds per hour

Qualifiers:

J - Result is estimated

U - below reported quantitation limit

_J - Laboratory data qualifier

_U - Data validation qualifier

System	Date	Temp (F)	Flow (scfm)
Off-Site	11/02/06	68	1370
SBPA	11/02/06	96	1693

Table 3.9
SBPA and Off-Site ISVE System Results
for Method TO-14 (VOCs) - December 2006
American Chemical Service
Griffith, Indiana

Compounds	Units	Sampled 12/11/06			
		SBPA ISVE		Off-Site ISVE	
1,1,1-Trichloroethane	ppbv	17,000		21,000	
1,1,2,2-Tetrachloroethane	ppbv	ND	U	ND	U
1,1,2-Trichloroethane	ppbv	66	J/J	160	J/J
1,1-Dichloroethane	ppbv	2,900		2,700	
1,1-Dichloroethene	ppbv	1,100		360	
1,2-Dichloroethane	ppbv	250		660	
1,2-Dichloropropane	ppbv	290		180	J/J
2-Butanone (Methyl Ethyl Ketone)	ppbv	1,000		16,000	
2-Hexanone	ppbv	ND	U	220	J/J
4-Methyl-2-pentanone	ppbv	1,600		7,100	
Acetone	ppbv	1,200		16,000	
Benzene	ppbv	5,200		12,000	
Eromodichloromethane	ppbv	ND	U	ND	U
Eromoform	ppbv	ND	U	ND	U
Eromomethane	ppbv	ND	U	ND	U
Carbon Disulfide	ppbv	77	J/J	440	J/J
Carbon Tetrachloride	ppbv	ND	U	ND	U
Chlorobenzene	ppbv	57	J/J	ND	U
Chloroethane	ppbv	640	/J	ND	U
Chloroform	ppbv	4,600		1,800	
Chloromethane	ppbv	ND	U	ND	U
cis-1,2-Dichloroethene	ppbv	14,000		1,400	
cis-1,3-Dichloropropene	ppbv	ND	U	ND	U
Dibromochloromethane	ppbv	ND	U	ND	U
Ethyl Benzene	ppbv	10,000		10,000	
m,p-Xylene	ppbv	44,000		44,000	
Methylene Chloride	ppbv	6,400		19,000	
c-Xylene	ppbv	20,000		17,000	
Styrene	ppbv	ND	U	ND	U
Tetrachloroethene	ppbv	25,000		16,000	
Toluene	ppbv	42,000		86,000	
trans-1,2-Dichloroethene	ppbv	120	J/J	ND	U
trans-1,3-Dichloropropene	ppbv	ND	U	ND	U
Trichloroethene	ppbv	13,000		13,000	
Vinyl Chloride	ppbv	1,300		ND	U
Total	ppbv	211,800		285,020	
Total	lb/hr	5.435		4.230	

Notes:

ND - Non-detect

ppbv - parts per billion volume

lb/hr - pounds per hour

Qualifiers:

J - Result is estimated

U - below reported quantitation limit

/J - Laboratory data qualifier

/ - Data validation qualifier

System	Date	Temp (F)	Flow (scfm)
Off-Site	12/11/06	64	1090
SBPA	12/11/06	98	1707

Table 3.10
Thermal Oxidizer 1 Results for Method TO-13 (SVOCs) - October 2006
American Chemical Service
Griffith, Indiana

Compounds	Units	Sampled 10/19/06						
		Therm-Ox 1				Destruction Efficiency		
		Influent	Influent Dup	Effluent		Low	High	Average
1,2,4-Trichlorobenzene	µg	0.72	J	0.91	J	ND	U	NC NC NC
1,2-Dichlorobenzene	µg	24		28		ND	U	100.00% 100.00% 100.00%
1,3-Dichlorobenzene	µg	2.6		2.9		ND	U	100.00% 100.00% 100.00%
1,4-Dichlorobenzene	µg	5.7		6.7		ND	U	100.00% 100.00% 100.00%
2,4,5-Trichlorophenol	µg	ND	U	ND	U	ND	U	NC NC NC
2,4,6-Trichlorophenol	µg	ND	U	ND	U	ND	U	NC NC NC
2,4-Dichlorophenol	µg	ND	U	ND	U	ND	U	NC NC NC
2,4-Dimethylphenol	µg	ND	U	ND	U	ND	U	NC NC NC
2,4-Dinitrophenol	µg	ND	U	ND	U	ND	U	NC NC NC
2,4-Dinitrotoluene	µg	ND	U	ND	U	ND	U	NC NC NC
2,6-Dinitrotoluene	µg	ND	U	ND	U	ND	U	NC NC NC
2-Chloronaphthalene	µg	ND	U	ND	U	ND	U	NC NC NC
2-Chlorophenol	µg	ND	U	ND	U	ND	U	NC NC NC
2-Methylnaphthalene	µg	14		19		ND	U	100.00% 100.00% 100.00%
2-Methylphenol (o-Cresol)	µg	ND	U	ND	U	ND	U	NC NC NC
2-Nitroaniline	µg	ND	U	ND	U	ND	U	NC NC NC
2-Nitrophenol	µg	ND	U	ND	U	ND	U	NC NC NC
3,3'-Dichlorobenzidine	µg	ND	U	ND	U	ND	U	NC NC NC
3-Nitroaniline	µg	ND	U	ND	U	ND	U	NC NC NC
4,6-Dinitro-2-methylphenol	µg	ND	U	ND	U	ND	U	NC NC NC
4-Bromophenyl-phenyl Ether	µg	ND	U	ND	U	ND	U	NC NC NC
4-Chloro-3-methylphenol	µg	ND	U	ND	U	ND	U	NC NC NC
4-Chloroaniline	µg	ND	U	ND	U	ND	U	NC NC NC
4-Chlorophenyl-phenyl Ether	µg	ND	U	ND	U	ND	U	NC NC NC
4-Methylphenol/3-Methylphenol	µg	ND	U	ND	U	ND	U	NC NC NC
4-Nitroaniline	µg	ND	U	ND	U	ND	U	NC NC NC
4-Nitrophenol	µg	ND	U	ND	U	ND	U	NC NC NC
Acenaphthene	µg	ND	U	ND	U	ND	U	NC NC NC
Acenaphthylene	µg	ND	U	ND	U	ND	U	NC NC NC
Anthracene	µg	ND	U	ND	U	ND	U	NC NC NC
Benz(a)anthracene	µg	ND	U	ND	U	ND	U	NC NC NC
Benz(a)pyrene	µg	ND	U	ND	U	ND	U	NC NC NC
Benz(b)fluoranthene	µg	ND	U	ND	U	ND	U	NC NC NC
Benz(g,h,i)perylene	µg	ND	U	ND	U	ND	U	NC NC NC
Benz(k)fluoranthene	µg	ND	U	ND	U	ND	U	NC NC NC
bis(2-Chloroethoxy) Methane	µg	ND	U	ND	U	ND	U	NC NC NC
bis(2-Chloroethyl) Ether	µg	ND		2.8		ND	U	NC NC NC
bis(2-Ethylhexyl)phthalate	µg	ND	U	ND	U	ND	U	NC NC NC
Butylbenzylphthalate	µg	ND	U	ND	U	ND	U	NC NC NC
Chrysene	µg	ND	U	ND	U	ND	U	NC NC NC
Dibenz(a,h)anthracene	µg	ND	U	ND	U	ND	U	NC NC NC
Dibenzofuran	µg	ND	U	ND	U	ND	U	NC NC NC
Diethylphthalate	µg	ND	U	I	J	0.93	J	NC NC NC
Dimethylphthalate	µg	ND	U	ND	U	ND	U	NC NC NC
di-n-Butylphthalate	µg	ND	U	I	J	ND	U	NC NC NC
Di-n-Octylphthalate	µg	ND	U	ND	U	ND	U	NC NC NC
Fluoranthene	µg	ND	U	ND	U	ND	U	NC NC NC
Fluorene	µg	ND	U	ND	U	ND	U	NC NC NC
Hexachlorobenzene	µg	ND	U	ND	U	ND	U	NC NC NC
Hexachlorobutadiene	µg	8.8		11		ND	U	100.00% 100.00% 100.00%
Hexachlorocyclopentadiene	µg	ND	U	ND	U	ND	U	NC NC NC
Hexachloroethane	µg	ND	U	ND	U	ND	U	NC NC NC
Indeno(1,2,3-c,d)pyrene	µg	ND	U	ND	U	ND	U	NC NC NC
Isochorone	µg	3		3.8		ND	U	100.00% 100.00% 100.00%

Table 3.10
Thermal Oxidizer 1 Results for Method TO-13 (SVOCs) - October 2006
American Chemical Service
Griffith, Indiana

Compounds	Units	Sampled 10/19/06							
		Therm-Ox 1				Effluent		Destruction Efficiency	
		Influent	Influent Dup	Low	High	Average			
Naphthalene	µg	25	32	ND	U	100.00%	100.00%	100.00%	
Nitrobenzene	µg	ND	U	ND	U	NC	NC	NC	
N-Nitroso-di-n-propylamine	µg	ND	U	ND	U	NC	NC	NC	
N-Nitrosodiphenylamine	µg	ND	U	ND	U	NC	NC	NC	
Pentachlorophenol	µg	ND	U	ND	U	NC	NC	NC	
Phenanthrene	µg	ND	U	ND	U	NC	NC	NC	
Phenol	µg	ND	U	ND	U	NC	NC	NC	
Pyrene	µg	ND	U	ND	U	NC	NC	NC	
Total	µg	83.82	109.11	0.93		98.89%	99.15%	99.02%	

Notes:

µg - Microgram

NC - Not calculated

ND - Non-detect

Qualifiers:

J - Result is estimated

U - below reported quantitation limit

/ - Laboratory data qualifier

/ - Data validation qualifier

Table 3.11
Thermal Oxidizer 1 Results for Method TO-13 (SVOCs) - November 2006
American Chemical Service
Griffith, Indiana

Compounds	Units	Sampled 11/2/06						
		Therm-Ox 1			Effluent		Low	High
		Influent	Influent Dup	Therm-Ox 1	Effluent	Effluent		
1,2,4-Trichlorobenzene	µg	NS		12	ND	U	NC	NC
1,2-Dichlorobenzene	µg	NS		26	ND	U	NC	NC
1,3-Dichlorobenzene	µg	NS		3.1	ND	U	NC	NC
1,4-Dichlorobenzene	µg	NS		6.6	ND	U	NC	NC
2,4,5-Trichlorophenol	µg	NS		ND	U	ND	NC	NC
2,4,6-Trichlorophenol	µg	NS		ND	U	ND	NC	NC
2,4-Dichlorophenol	µg	NS		ND	U	ND	U	NC
2,4-Dimethylphenol	µg	NS		ND	U	ND	U	NC
2,4-Dinitrophenol	µg	NS		ND	U	ND	U	NC
2,4-Dinitrotoluene	µg	NS		ND	U	ND	U	NC
2,6-Dinitrotoluene	µg	NS		ND	U	ND	U	NC
2-Chloronaphthalene	µg	NS		ND	U	ND	U	NC
2-Chlorophenol	µg	NS		ND	U	ND	U	NC
2-Methylnaphthalene	µg	NS		9.6		ND	U	NC
2-Methylphenol (o-Cresol)	µg	NS		ND	U	ND	U	NC
2-Nitroaniline	µg	NS		ND	U	ND	U	NC
2-Nitrophenol	µg	NS		ND	U	ND	U	NC
3,3'-Dichlorobenzidine	µg	NS		ND	U	ND	U	NC
3-Nitroaniline	µg	NS		ND	U	ND	U	NC
4,6-Dinitro-2-methylphenol	µg	NS		ND	U	ND	U	NC
4-Eromoethyl-phenyl Ether	µg	NS		ND	U	ND	U	NC
4-Chloro-3-methylphenol	µg	NS		ND	U	ND	U	NC
4-Chloroaniline	µg	NS		ND	U	ND	U	NC
4-Chlorophenyl-phenyl Ether	µg	NS		ND	U	ND	U	NC
4-Methylphenol/3-Methylphenol	µg	NS		ND	U	ND	U	NC
4-Nitroaniline	µg	NS		ND	U	ND	U	NC
4-Nitrophenol	µg	NS		ND	U	ND	U	NC
Acenaphthene	µg	NS		ND	U	ND	U	NC
Acenaphthylene	µg	NS		ND	U	ND	U	NC
Anthracene	µg	NS		ND	U	ND	U	NC
Benzo(a)anthracene	µg	NS		ND	U	ND	U	NC
Benzo(a)pyrene	µg	NS		ND	U	ND	U	NC
Benzo(b)fluoranthene	µg	NS		ND	U	ND	U	NC
Benzo(g,h,i)perylene	µg	NS		ND	U	ND	U	NC
Benzo(k)fluoranthene	µg	NS		ND	U	ND	U	NC
bis(2-Chloroethoxy) Methane	µg	NS		ND	U	ND	U	NC
bis(2-Chloroethyl) Ether	µg	NS		2.9		ND	U	NC
bis(2-Ethylhexyl)phthalate	µg	NS		1.5	J/J	ND	U	NC
Butylbenzylphthalate	µg	NS		ND	U	ND	U	NC
Chrysene	µg	NS		ND	U	ND	U	NC
Di- <i>benz(a,h)</i> anthracene	µg	NS		ND	U	ND	U	NC
Di- <i>benzofuran</i>	µg	NS		ND	U	ND	U	NC
Diethylphthalate	µg	NS		ND	U	ND	J/J	NC
Dimethylphthalate	µg	NS		ND	U	ND	U	NC
di-n-Butylphthalate	µg	NS		1.1	J/J	1.3	J/J	NC
Di-n-Octylphthalate	µg	NS		ND	U	ND	U	NC
Fluoranthene	µg	NS		ND	U	ND	U	NC
Fluorene	µg	NS		ND	U	ND	U	NC
Hexachlorobenzene	µg	NS		ND	U	ND	U	NC
Hexachlorobutadiene	µg	NS		11		ND	U	NC
Hexachlorocyclopentadiene	µg	NS		ND	U	ND	U	NC
Hexachloroethane	µg	NS		ND	U	ND	U	NC
Indeno(1,2,3-c,d)pyrene	µg	NS		ND	U	ND	U	NC
Isophorone	µg	NS		3		ND	U	NC

Table 3.11
Thermal Oxidizer 1 Results for Method TO-13 (SVOCs) - November 2006
American Chemical Service
Griffith, Indiana

Compounds	Units	Sampled 11/2/06						
		Therm-Ox 1				Destruction Efficiency		
		Influent	Influent Dup	Effluent		Low	High	Average
Naphthalene	µg	NS	18	ND	U	NC	NC	NC
Nitrobenzene	µg	NS	ND	U	ND	U	NC	NC
N-Nitroso-di-n-propylamine	µg	NS	ND	U	ND	U	NC	NC
N-Nitrosodiphenylamine	µg	NS	ND	U	ND	U	NC	NC
Pentachlorophenol	µg	NS	ND	U	ND	U	NC	NC
Phenanthrene	µg	NS	ND	U	ND	U	NC	NC
Phenol	µg	NS	ND	U	ND	U	NC	NC
Pyrene	µg	NS	ND	U	ND	U	NC	NC
Total	µg	NS	94.80		1.3		NC	NC

Notes:

µg - Microgram

NC - Not calculated

ND - Non-detect

Qualifiers:

J - Result is estimated

U - below reported quantitation limit

/ - Laboratory data qualifier

/ - Data validation qualifier

Table 3.12
Thermal Oxidizer 1 Results for Method TO-13 (SVOCs) - December 2006
American Chemical Service
Griffith, Indiana

Compounds	Units	Sampled 12/11/06							
		Therm-Ox 1				Destruction Efficiency			
		Influent	Influent Dup	Effluent		Low	High	Average	
1,2,4-Trichlorobenzene	µg	13		15		ND	U	100.00%	100.00%
1,2-Dichlorobenzene	µg	23		25		ND	U	100.00%	100.00%
1,3-Dichlorobenzene	µg	2.8		3		ND	U	100.00%	100.00%
1,4-Dichlorobenzene	µg	5.8		6		ND	U	100.00%	100.00%
2,4,5-Trichlorophenol	µg	ND	U	ND	U	ND	U	NC	NC
2,4,6-Trichlorophenol	µg	ND	U	ND	U	ND	U	NC	NC
2,4-Dichlorophenol	µg	ND	U	ND	U	ND	U	NC	NC
2,4-Dimethylphenol	µg	ND	U	ND	U	ND	U	NC	NC
2,4-Dinitrophenol	µg	ND	U	ND	U	ND	U	NC	NC
2,4-Dinitrotoluene	µg	ND	U	ND	U	ND	U	NC	NC
2,6-Dinitrotoluene	µg	ND	U	ND	U	ND	U	NC	NC
2-Chloronaphthalene	µg	ND	U	ND	U	ND	U	NC	NC
2-Chlorophenol	µg	ND	U	ND	U	ND	U	NC	NC
2-Methylnaphthalene	µg	10		11		ND	U	100.00%	100.00%
2-Methylphenol (o-Cresol)	µg	ND	U	ND	U	ND	U	NC	NC
2-Nitroaniline	µg	ND	U	ND	U	ND	U	NC	NC
2-Nitrophenol	µg	ND	U	ND	U	ND	U	NC	NC
3,3-Dichlorobenzidine	µg	ND	U	ND	U	ND	U	NC	NC
3-Nitroaniline	µg	ND	U	ND	U	ND	U	NC	NC
4,6-Dinitro-2-methylphenol	µg	ND	U	ND	U	ND	U	NC	NC
4-Bromophenyl-phenyl Ether	µg	ND	U	ND	U	ND	U	NC	NC
4-Chloro-3-methylphenol	µg	ND	U	ND	U	ND	U	NC	NC
4-Chloroaniline	µg	ND	U	ND	U	ND	U	NC	NC
4-Chlorophenyl-phenyl Ether	µg	ND	U	ND	U	ND	U	NC	NC
4-Methylphenol/3-Methylphenol	µg	ND	U	ND	U	ND	U	NC	NC
4-Nitroaniline	µg	ND	U	ND	U	ND	U	NC	NC
4-Nitrophenol	µg	ND	U	ND	U	ND	U	NC	NC
Acenaphthene	µg	ND	U	ND	U	ND	U	NC	NC
Acenaphthylene	µg	ND	U	ND	U	ND	U	NC	NC
Anthracene	µg	ND	U	ND	U	ND	U	NC	NC
Benzo(a)anthracene	µg	ND	U	ND	U	ND	U	NC	NC
Benzo(a)pyrene	µg	ND	U	ND	U	ND	U	NC	NC
Benzo(b)fluoranthene	µg	ND	U	ND	U	ND	U	NC	NC
Benzo(g,h,i)perylene	µg	ND	U	ND	U	ND	U	NC	NC
Benzo(k)fluoranthene	µg	ND	U	ND	U	ND	U	NC	NC
bis(2-Chloroethoxy) Methane	µg	ND	U	ND	U	ND	U	NC	NC
bis(2-Chloroethyl) Ether	µg	2.6		2.7		ND	U	100.00%	100.00%
bis(2-Ethylhexyl)phthalate	µg	ND	U	ND	U	ND	U	NC	NC
Buylbenzylphthalate	µg	ND	U	ND	U	ND	U	NC	NC
Chrysene	µg	ND	U	ND	U	ND	U	NC	NC
Dibenzo(a,h)anthracene	µg	ND	U	ND	U	ND	U	NC	NC
Dibenzofuran	µg	ND	U	ND	U	ND	U	NC	NC
Diethylphthalate	µg	ND	U	ND	U	ND	U	NC	NC
Dimethylphthalate	µg	ND	U	ND	U	ND	U	NC	NC
di-1-Butylphthalate	µg	ND	U	ND	U	ND	U	NC	NC
Di-n-Octylphthalate	µg	ND	U	ND	U	ND	U	NC	NC
Fluoranthene	µg	ND	U	ND	U	ND	U	NC	NC
Fluorene	µg	ND	U	ND	U	ND	U	NC	NC
Hexachlorobenzene	µg	ND	U	ND	U	ND	U	NC	NC
Hexachlorobutadiene	µg	10		11		ND	U	100.00%	100.00%
Hexachlorocyclopentadiene	µg	ND	U	ND	U	ND	U	NC	NC
Hexachloroethane	µg	ND	U	ND	U	ND	U	NC	NC
Indeno(1,2,3-c,d)pyrene	µg	ND	U	ND	U	ND	U	NC	NC
Isophorone	µg	2.4		2.4		ND	U	100.00%	100.00%

Table 3.12
Thermal Oxidizer 1 Results for Method TO-13 (SVOCs) - December 2006
American Chemical Service
Griffith, Indiana

Compounds	Units	Sampled 12/11/06							
		Therm-Ox 1				Destruction Efficiency			
		Influent	Influent Dup	Effluent		Low	High	Average	
Naphthalene	µg	17	19	ND	U	100.00%	100.00%	100.00%	
Nitrobenzene	µg	ND	U	ND	U	NC	NC	NC	
N-Nitroso-di-n-propylamine	µg	ND	U	ND	U	NC	NC	NC	
N-Nitrosodiphenylamine	µg	ND	U	ND	U	NC	NC	NC	
Pentachlorophenol	µg	ND	U	ND	U	NC	NC	NC	
Phenanthrene	µg	ND	U	ND	U	NC	NC	NC	
Phenol	µg	ND	U	ND	U	NC	NC	NC	
Pyrene	µg	ND	U	ND	U	NC	NC	NC	
Total	µg			95.10		0.0	NC	NC	NC

Notes:

µg - Microgram

NC - Not calculated

ND - Non-detect

Qualifiers:

U - below reported quantitation limit

/ - Laboratory data qualifier

/_ - Data validation qualifier

Table 3.13
Thermal Oxidizer 2 Results for Method TO-13 (SVOCs) - October 2006
American Chemical Service
Griffith, Indiana

Compounds	Units	Sampled 10/19/06					
		Therm-Ox 2			Destruction Efficiency		
		Influent	Influent Dup	Effluent	Low	High	Average
1,2,4-Trichlorobenzene	µg	1.2		1.7	ND	U	100.00% 100.00% 100.00%
1,2-Dichlorobenzene	µg	42		52	ND	U	100.00% 100.00% 100.00%
1,3-Dichlorobenzene	µg	1.4		1.7	ND	U	100.00% 100.00% 100.00%
1,4-Dichlorobenzene	µg	5.1		6.3	ND	U	100.00% 100.00% 100.00%
2,4,5-Trichlorophenol	µg	ND	U	ND	ND	U	NC NC NC
2,4,6-Trichlorophenol	µg	ND	U	ND	ND	U	NC NC NC
2,4-Dichlorophenol	µg	ND	U	ND	U	ND	NC NC NC
2,4-Dimethylphenol	µg	ND	U	ND	U	ND	NC NC NC
2,4-Dinitrophenol	µg	ND	U	ND	U	ND	NC NC NC
2,4-Dinitrotoluene	µg	ND	U	ND	U	ND	NC NC NC
2,6-Dinitrotoluene	µg	ND	U	ND	U	ND	NC NC NC
2-Chloronaphthalene	µg	ND	U	ND	U	ND	NC NC NC
2-Chlorophenol	µg	ND	U	ND	U	ND	NC NC NC
2-Methylnaphthalene	µg	4.8		6	ND	U	100.00% 100.00% 100.00%
2-Methylphenol (o-Cresol)	µg	ND	U	ND	U	ND	NC NC NC
2-Nitroaniline	µg	ND	U	ND	U	ND	NC NC NC
2-Nitrophenol	µg	ND	U	ND	U	ND	NC NC NC
3,3'-Dichlorobenzidine	µg	ND	U	ND	U	ND	NC NC NC
3-Nitroaniline	µg	ND	U	ND	U	ND	NC NC NC
4,6-Dinitro-2-methylphenol	µg	ND	U	ND	U	ND	NC NC NC
4-Bromophenyl-phenyl Ether	µg	ND	U	ND	U	ND	NC NC NC
4-Chloro-3-methylphenol	µg	ND	U	ND	U	ND	NC NC NC
4-Chloroaniline	µg	ND	U	ND	U	ND	NC NC NC
4-Chlorophenyl-phenyl Ether	µg	ND	U	ND	U	ND	NC NC NC
4-Methylphenol/3-Methylphenol	µg	ND	U	ND	U	ND	NC NC NC
4-Nitroaniline	µg	ND	U	ND	U	ND	NC NC NC
4-Nitrophenol	µg	ND	U	ND	U	ND	NC NC NC
Acenaphthene	µg	ND	U	ND	U	ND	NC NC NC
Acenaphthylene	µg	ND	U	ND	U	ND	NC NC NC
Anthracene	µg	ND	U	ND	U	ND	NC NC NC
Benz(a)anthracene	µg	ND	U	ND	U	ND	NC NC NC
Benz(a)pyrene	µg	ND	U	ND	U	ND	NC NC NC
Benz(b)fluoranthene	µg	ND	U	ND	U	ND	NC NC NC
Benz(g,h,i)perylene	µg	ND	U	ND	U	ND	NC NC NC
Benz(k)fluoranthene	µg	ND	U	ND	U	ND	NC NC NC
bis(2-Chloroethoxy) Methane	µg	ND	U	ND	U	ND	NC NC NC
bis(2-Chloroethyl) Ether	µg	3.4		4.4	ND	U	100.00% 100.00% 100.00%
bis(2-Ethylhexyl)phthalate	µg	12		2.6	J	1.4 J	NC NC NC NC
Butylbenzylphthalate	µg	ND	U	ND	U	ND	NC NC NC NC
Chrysene	µg	ND	U	ND	U	ND	NC NC NC NC
Dibenz(a,h)anthracene	µg	ND	U	ND	U	ND	NC NC NC NC
Dibenzofuran	µg	ND	U	ND	U	ND	NC NC NC NC
Diethylphthalate	µg	0.81	J	1.1 J	0.92	J	NC NC NC NC
Dimethylphthalate	µg	ND	U	ND	U	ND	NC NC NC NC
di-n-Butylphthalate	µg	ND	U	0.89	J	ND U	NC NC NC NC
Di-n-Octylphthalate	µg	ND	U	ND	U	ND	NC NC NC NC
Fluoranthene	µg	ND	U	ND	U	ND	NC NC NC NC
Fluorene	µg	ND	U	ND	U	ND	NC NC NC NC
Hexachlorobenzene	µg	ND	U	ND	U	ND	NC NC NC NC
Hexachlorobutadiene	µg	2.1		2.9	ND	U	100.00% 100.00% 100.00%
Hexachlorocyclopentadiene	µg	ND	U	ND	U	ND	NC NC NC NC
Hexachloroethane	µg	ND	U	ND	U	ND	NC NC NC NC
Indeno(1,2,3-c,d)pyrene	µg	ND	U	ND	U	ND	NC NC NC NC
Isophorone	µg	12		16	ND	U	100.00% 100.00% 100.00%

Table 3.13
Thermal Oxidizer 2 Results for Method TO-13 (SVOCs) - October 2006
American Chemical Service
Griffith, Indiana

Compounds	Units	Sampled 10/19/06								
		Therm-Ox 2				Destruction Efficiency				
		Influent		Influent Dup		Effluent		Low	High	Average
Naphthalene	µg	24		32		ND	U	100.00%	100.00%	100.00%
Nitrobenzene	µg	ND	U	ND	U	ND	U	NC	NC	NC
N-Nitroso-di-n-propylamine	µg	ND	U	ND	U	ND	U	NC	NC	NC
N-Nitrosodiphenylamine	µg	ND	U	ND	U	ND	U	NC	NC	NC
Pentachlorophenol	µg	ND	U	ND	U	ND	U	NC	NC	NC
Phenanthrene	µg	ND	U	ND	U	ND	U	NC	NC	NC
Phenol	µg	ND	U	ND	U	ND	U	NC	NC	NC
Pyrene	µg	ND	U	ND	U	ND	U	NC	NC	NC
Total	µg	108.81		127.59		2.32		97.87%	98.18%	98.02%

Notes:

µg - Microgram

NC - Not calculated

ND - Non-detect

Qualifiers:

J - Result is estimated

U - below reported quantitation limit

/ - Laboratory data qualifier

/ - Data validation qualifier

Table 3.14
Thermal Oxidizer 2 Results for Method TO-13 (SVOCs) - November 2006
American Chemical Service
Griffith, Indiana

Compounds	Units	Sampled 11/2/06						
		Therm-Ox 2			Effluent		Low	High
		Influent	Influent Dup	J/J	ND	U		
1,2,4-Trichlorobenzene	µg	NS	0.67	J/J	ND	U	NC	NC
1,2-Dichlorobenzene	µg	NS	24		1.4		NC	NC
1,3-Dichlorobenzene	µg	NS	1		ND	U	NC	NC
1,4-Dichlorobenzene	µg	NS	3.2		ND	U	NC	NC
2,4,5-Trichlorophenol	µg	NS	ND	U	ND	U	NC	NC
2,4,6-Trichlorophenol	µg	NS	ND	U	ND	U	NC	NC
2,4-Dichlorophenol	µg	NS	ND	U	ND	U	NC	NC
2,4-Dimethylphenol	µg	NS	ND	U	ND	U	NC	NC
2,4-Dinitrophenol	µg	NS	ND	U	ND	U	NC	NC
2,4-Dinitrotoluene	µg	NS	ND	U	ND	U	NC	NC
2-Chloronaphthalene	µg	NS	ND	U	ND	U	NC	NC
2-Chlorophenol	µg	NS	ND	U	ND	U	NC	NC
2-Methylnaphthalene	µg	NS	2.2		ND	U	NC	NC
2-Methylphenol (o-Cresol)	µg	NS	ND	U	ND	U	NC	NC
2-Nitroaniline	µg	NS	ND	U	ND	U	NC	NC
2-Nitrophenol	µg	NS	ND	U	ND	U	NC	NC
3,3'-Dichlorobenzidine	µg	NS	ND	U	ND	U	NC	NC
3-Nitroaniline	µg	NS	ND	U	ND	U	NC	NC
4,6-Dinitro-2-methylphenol	µg	NS	ND	U	ND	U	NC	NC
4-Bromophenyl-phenyl Ether	µg	NS	ND	U	ND	U	NC	NC
4-Chloro-3-methylphenol	µg	NS	ND	U	ND	U	NC	NC
4-Chloroaniline	µg	NS	ND	U	ND	U	NC	NC
4-Chlorophenyl-phenyl Ether	µg	NS	ND	U	ND	U	NC	NC
4-Methylphenol/3-Methylphenol	µg	NS	ND	U	ND	U	NC	NC
4-Nitroaniline	µg	NS	ND	U	ND	U	NC	NC
4-Nitrophenol	µg	NS	ND	U	ND	U	NC	NC
Acenaphthene	µg	NS	ND	U	ND	U	NC	NC
Acenaphthylene	µg	NS	ND	U	ND	U	NC	NC
Anthracene	µg	NS	ND	U	ND	U	NC	NC
Benz(a)anthracene	µg	NS	ND	U	ND	U	NC	NC
Benz(a)pyrene	µg	NS	ND	U	ND	U	NC	NC
Benz(b)fluoranthene	µg	NS	ND	U	ND	U	NC	NC
Benz(g,h,i)perylene	µg	NS	ND	U	ND	U	NC	NC
Benz(k)fluoranthene	µg	NS	ND	U	ND	U	NC	NC
bis(2-Chloroethoxy) Methane	µg	NS	ND	U	ND	U	NC	NC
bis(2-Chloroethyl) Ether	µg	NS	2.2		ND	U	NC	NC
bis(2-Ethylhexyl)phthalate	µg	NS	1.1	J/J	3.8	J/J	NC	NC
Butylbenzylphthalate	µg	NS	ND	U	ND	U	NC	NC
Chrysene	µg	NS	ND	U	ND	U	NC	NC
Dibenz(a,h)anthracene	µg	NS	ND	U	ND	U	NC	NC
Dibenzofuran	µg	NS	ND	U	ND	U	NC	NC
Diethylphthalate	µg	NS	ND	U	ND	U	NC	NC
Dimethylphthalate	µg	NS	ND	U	ND	U	NC	NC
di- <i>t</i> -Butylphthalate	µg	NS	1.1	J/J	1.1	J/J	NC	NC
Di- <i>t</i> -Octylphthalate	µg	NS	ND	U	ND	U	NC	NC
Fluoranthene	µg	NS	ND	U	ND	U	NC	NC
Fluorene	µg	NS	ND	U	ND	U	NC	NC
Heptachlorobenzene	µg	NS	ND	U	ND	U	NC	NC
Heptachlorobutadiene	µg	NS	1.2		ND	U	NC	NC
Heptachlorocyclopentadiene	µg	NS	ND	U	ND	U	NC	NC
Heptachloroethane	µg	NS	ND	U	ND	U	NC	NC
Indeno(1,2,3-c,d)pyrene	µg	NS	ND	U	ND	U	NC	NC
Isophorone	µg	NS	5.3		ND	U	NC	NC

Table 3.14
Thermal Oxidizer 2 Results for Method TO-13 (SVOCs) - November 2006
American Chemical Service
Griffith, Indiana

		Sampled 11/2/06							
		Therm-Ox 2				Destruction Efficiency			
Compounds	Units	Influent	Influent Dup	Effluent		Low	High	Average	
Naphthalene	µg	NS	12	2		NC	NC	NC	
Nitrobenzene	µg	NS	ND	U	ND	U	NC	NC	NC
N-Nitroso-di-n-propylamine	µg	NS	ND	U	ND	U	NC	NC	NC
N-Nitrosodiphenylamine	µg	NS	ND	U	ND	U	NC	NC	NC
Pentachlorophenol	µg	NS	ND	U	ND	U	NC	NC	NC
Phenanthrene	µg	NS	ND	U	ND	U	NC	NC	NC
Phenol	µg	NS	ND	U	ND	U	NC	NC	NC
Pyrene	µg	NS	ND	U	ND	U	NC	NC	NC
Total	µg	0.00	53.97		8.30		NC	NC	NC

Note:

µg - Microgram
 NC - Not calculated
 ND - Non-detect

Qualifiers:

J - Result is estimated
 U - below reported quantitation limit
 J - Laboratory data qualifier
 U - Data validation qualifier

Table 3.15
Thermal Oxidizer 2 Results for Method TO-13 (SVOCs) - December 2006
American Chemical Service
Griffith, Indiana

Compounds	Units	Sampled 12/11/06					
		Therm-Ox 2			Destruction Efficiency		
		Influent	Influent Dup	Effluent	Low	High	Average
1,2,4-Trichlorobenzene	µg	1.1		2.4	ND	U	100.00%
1,2-Dichlorobenzene	µg	39		72	3.2	U	91.79%
1,3-Dichlorobenzene	µg	1.6		3	ND	U	100.00%
1,4-Dichlorobenzene	µg	5.1		9.1	ND	U	100.00%
2,4,5-Trichlorophenol	µg	ND	U	ND	U	ND	NC
2,4,6-Trichlorophenol	µg	ND	U	ND	U	ND	NC
2,4-Dichlorophenol	µg	ND	U	ND	U	ND	NC
2,4-Dimethylphenol	µg	ND	U	ND	U	ND	NC
2,4-Dinitrophenol	µg	ND	U	ND	U	ND	NC
2,4-Dinitrotoluene	µg	ND	U	ND	U	ND	NC
2,6-Dinitrotoluene	µg	ND	U	ND	U	ND	NC
2-Chloronaphthalene	µg	ND	U	ND	U	ND	NC
2-Chlorophenol	µg	ND	U	ND	U	ND	NC
2-Methylnaphthalene	µg	4.1		8	0.65	J/J	NC
2-Methylphenol (o-Cresol)	µg	ND	U	ND	U	ND	NC
2-Nitroaniline	µg	ND	U	ND	U	ND	NC
2-Nitrophenol	µg	ND	U	ND	U	ND	NC
3,3'-Dichlorobenzidine	µg	ND	U	ND	U	ND	NC
3-Nitroaniline	µg	ND	U	ND	U	ND	NC
4,6-Dinitro-2-methylphenol	µg	ND	U	ND	U	ND	NC
4-Bromophenyl-phenyl Ether	µg	ND	U	ND	U	ND	NC
4-Chloro-3-methylphenol	µg	ND	U	ND	U	ND	NC
4-Chloroaniline	µg	ND	U	ND	U	ND	NC
4-Chlorophenyl-phenyl Ether	µg	ND	U	ND	U	ND	NC
4-Methylphenol/3-Methylphenol	µg	ND	U	ND	U	ND	NC
4-Nitroaniline	µg	ND	U	ND	U	ND	NC
4-Nitrophenol	µg	ND	U	ND	U	ND	NC
Acenaphthene	µg	ND	U	ND	U	ND	NC
Acenaphthylene	µg	ND	U	ND	U	ND	NC
Anthracene	µg	ND	U	ND	U	ND	NC
Benzo(a)anthracene	µg	ND	U	ND	U	ND	NC
Benzo(a)pyrene	µg	ND	U	ND	U	ND	NC
Benzo(b)fluoranthene	µg	ND	U	ND	U	ND	NC
Benzo(g,h,i)perylene	µg	ND	U	ND	U	ND	NC
Benzo(k)fluoranthene	µg	ND	U	ND	U	ND	NC
bis(2-Chloroethyl) Methane	µg	ND	U	ND	U	ND	NC
bis(2-Chloroethyl) Ether	µg	3.1		6.1	ND	U	100.00%
bis(2-Ethylhexyl)phthalate	µg	1.3	J/J	1.4	J/J	7.8	NC
Butylbenzylphthalate	µg	ND	U	ND	U	ND	NC
Chrysene	µg	ND	U	ND	U	ND	NC
Dibenz(a,h)anthracene	µg	ND	U	ND	U	ND	NC
Dibenzo furan	µg	ND	U	ND	U	ND	NC
Diethylphthalate	µg	ND	U	ND	U	ND	NC
Dimethylphthalate	µg	ND	U	ND	U	ND	NC
di-n-Butylphthalate	µg	ND	U	ND	U	ND	NC
Di-n-Octylphthalate	µg	ND	U	ND	U	ND	NC
Fluoranthene	µg	ND	U	ND	U	ND	NC
Fluorene	µg	ND	U	ND	U	ND	NC
Hexachlorobenzene	µg	ND	U	ND	U	ND	NC
Hexachlorobutadiene	µg	2.1		4.4	ND	U	100.00%
Hexachlorocyclopentadiene	µg	ND	U	ND	U	ND	NC
Hexachloroethane	µg	ND	U	ND	U	ND	NC
Indeno(1,2,3-c,d)pyrene	µg	ND	U	ND	U	ND	NC
Isophorone	µg	11		22	ND	U	100.00%

Table 3.15
Thermal Oxidizer 2 Results for Method TO-13 (SVOCs) - December 2006
American Chemical Service
Griffith, Indiana

Compounds	Units	Sampled 12/11/06						
		Therm-Ox 2				Destruction Efficiency		
		Influent	Influent Dup	Effluent		Low	High	Average
Naphthalene	µg	22		4.6		79.09%	89.55%	84.32%
Nitrobenzene	µg	2.7		ND	U	ND	U	NC
N-Nitroso-di-n-propylamine	µg	ND	U	ND	U	ND	U	NC
N-Nitrosodiphenylamine	µg	ND	U	ND	U	ND	U	NC
Pentachlorophenol	µg	ND	U	ND	U	ND	U	NC
Phenanthrene	µg	ND	U	ND	U	ND	U	NC
Phenol	µg	ND	U	3.2	J/J	ND	U	NC
Pyrene	µg	ND	U	ND	U	ND	U	NC
Total	µg	93.10		175.60		16.25		82.55% 90.75% 86.65%

Notes:

µg - Microgram

NC - Not calculated

ND - Non-detect

Qualifiers:

J - Result is estimated

U - below reported quantitation limit

J - Laboratory data qualifier

/ - Data validation qualifier

Table 3.16
SBPA and Off-Site ISVE System Results
for Method TO-13 (SVOCs) - October 2006
American Chemical Service
Griffith, Indiana

Compounds	Units	Sampled 10/19/2006			
		SBPA ISVE		Off-Site ISVE	
1,2,4-Trichlorobenzene	µg	0.53	J	4.5	
1,2-Dichlorobenzene	µg	22		88	
1,3-Dichlorobenzene	µg	2.3		2.6	
1,4-Dichlorobenzene	µg	5		10	
2,4,5-Trichlorophenol	µg	ND	U	ND	U
2,4,6-Trichlorophenol	µg	ND	U	ND	U
2,4-Dichlorophenol	µg	ND	U	ND	U
2,4-Dimethylphenol	µg	ND	U	ND	U
2,4-Dinitrophenol	µg	ND	U	ND	U
2,4-Dinitrotoluene	µg	ND	U	ND	U
2,6-Dinitrotoluene	µg	ND	U	ND	U
2-Chloronaphthalene	µg	ND	U	ND	U
2-Chlorophenol	µg	ND	U	ND	U
2-Methylnaphthalene	µg	12		21	
2-Methylphenol (o-Cresol)	µg	ND	U	ND	U
2-Nitroaniline	µg	ND	U	ND	U
2-Nitrophenol	µg	ND	U	ND	U
3,3'-Dichlorobenzidine	µg	ND	U	ND	U
3-Nitroaniline	µg	ND	U	ND	U
4,6-Dinitro-2-methylphenol	µg	ND	U	ND	U
4-Bromophenyl-phenyl Ether	µg	ND	U	ND	U
4-Chloro-3-methylphenol	µg	ND	U	ND	U
4-Chloroaniline	µg	ND	U	ND	U
4-Chlorophenyl-phenyl Ether	µg	ND	U	ND	U
4-Methylphenol/3-Methylphenol	µg	ND	U	4.8	J
4-Nitroaniline	µg	ND	U	ND	U
4-Nitrophenol	µg	ND	U	ND	U
Acenaphthene	µg	ND	U	ND	U
Acenaphthylene	µg	ND	U	ND	U
Anthracene	µg	ND	U	ND	U
Benz(a)anthracene	µg	ND	U	ND	U
Benz(a)pyrene	µg	ND	U	ND	U
Benz(b)fluoranthene	µg	ND	U	ND	U
Benz(g,h,i)perylene	µg	ND	U	ND	U
Benz(k)fluoranthene	µg	ND	U	ND	U
bis(2-Chloroethoxy) Methane	µg	ND	U	ND	U
bis(2-Chloroethyl) Ether	µg	2.1		9.2	
bis(2-Ethylhexyl)phthalate	µg	14		4.2	J
Butylbenzylphthalate	µg	ND	U	ND	U
Chrysene	µg	ND	U	ND	U
Dibenz(a,h)anthracene	µg	ND	U	ND	U
Dibenzofuran	µg	ND	U	ND	U
Diethylphthalate	µg	1.3	J	ND	U
Dimethylphthalate	µg	ND	U	ND	U
di-n-Butylphthalate	µg	ND	U	0.99	J
Di-n-Octylphthalate	µg	ND	U	ND	U
Fluoranthene	µg	ND	U	ND	U
Fluorene	µg	ND	U	ND	U
Hexachlorobenzene	µg	ND	U	ND	U
Hexachlorobutadiene	µg	8.2		5.9	
Hexachlorocyclopentadiene	µg	ND	U	ND	U
Hexachloroethane	µg	ND	U	ND	U
Indeno(1,2,3-c,d)pyrene	µg	ND	U	ND	U
Isophorone	µg	5.9		49	

Table 3.16
SBPA and Off-Site ISVE System Results
for Method TO-13 (SVOCs) - October 2006
American Chemical Service
Griffith, Indiana

Compounds	Units	Sampled 10/19/2006			
		SBPA ISVE	Off-Site ISVE		
Naphthalene	µg	22		100	
Nitrobenzene	µg	ND	U	ND	U
N-Nitroso-di-n-propylamine	µg	ND	U	ND	U
N-Nitrosodiphenylamine	µg	ND	U	ND	U
Pentachlorophenol	µg	ND	U	ND	U
Phenanthrene	µg	ND	U	ND	U
Phenol	µg	ND	U	ND	U
Pyrene	µg	ND	U	ND	U
Total	µg	95.33		300.19	

Notes:

µg - Microgram

NC - Not calculated

ND - Non-detect

Qualifiers:

J - Result is estimated

U - below reported quantitation limit

/ - Laboratory data qualifier

/_ - Data validation qualifier

Table 3.17
SBPA and Off-Site ISVE System Results
for Method TO-13 (SVOCs) - November 2006
American Chemical Service
Griffith, Indiana

Compounds	Units	Sampled 11/2/2006		
		SBPA ISVE	Off-Site ISVE	
1,2,4-Trichlorobenzene	µg	11		4.4
1,2-Dichlorobenzene	µg	24		77
1,3-Dichlorobenzene	µg	2.7		2.7
1,4-Dichlorobenzene	µg	5.7		10
2,4,5-Trichlorophenol	µg	ND	U	ND U
2,4,6-Trichlorophenol	µg	ND	U	ND U
2,4-Dichlorophenol	µg	ND	U	ND U
2,4-Dimethylphenol	µg	ND	U	ND U
2,4-Dinitrophenol	µg	ND	U	ND U
2,4-Dinitrotoluene	µg	ND	U	ND U
2,6-Dinitrotoluene	µg	ND	U	ND U
2-Chloronaphthalene	µg	ND	U	ND U
2-Chlorophenol	µg	ND	U	ND U
2-Methylnaphthalene	µg	9.1		20
2-Methylphenol (o-Cresol)	µg	ND	U	ND U
2-Nitroaniline	µg	ND	U	ND U
2-Nitrophenol	µg	ND	U	ND U
3,3'-Dichlorobenzidine	µg	ND	U	ND U
3-Nitroaniline	µg	ND	U	ND U
4,6-Dinitro-2-methylphenol	µg	ND	U	ND U
4-Bromophenyl-phenyl Ether	µg	ND	U	ND U
4-Chloro-3-methylphenol	µg	ND	U	ND U
4-Chloroaniline	µg	ND	U	ND U
4-Chlorophenyl-phenyl Ether	µg	ND	U	ND U
4-Methylphenol/3-Methylphenol	µg	ND	U	ND U
4-Nitroaniline	µg	ND	U	ND U
4-Nitrophenol	µg	ND	U	ND U
Acenaphthene	µg	ND	U	ND U
Acenaphthylene	µg	ND	U	ND U
Anthracene	µg	ND	U	ND U
Benzo(a)anthracene	µg	ND	U	ND U
Benzo(a)pyrene	µg	ND	U	ND U
Benzo(b)fluoranthene	µg	ND	U	ND U
Benzo(g,h,i)perylene	µg	ND	U	ND U
Benzo(k)fluoranthene	µg	ND	U	ND U
bis(2-Chloroethoxy) Methane	µg	ND	U	ND U
bis(2-Chloroethyl) Ether	µg	3.2		ND U
bis(2-Ethylhexyl)phthalate	µg	4.4	J/J	5.8
Butylbenzylphthalate	µg	ND	U	ND U
Chrysene	µg	ND	U	ND U
Dibenz(a,h)anthracene	µg	ND	U	ND U
Dibenzofuran	µg	ND	U	ND U
Diethylphthalate	µg	ND	U	ND U
Dimethylphthalate	µg	ND	U	ND U
di-n-Butylphthalate	µg	1.4	J/J	2.9 J/J
Di-n-Octylphthalate	µg	ND	U	ND U
Fluoranthene	µg	ND	U	ND U
Fluorene	µg	ND	U	ND U
Hexachlorobenzene	µg	ND	U	ND U
Hexachlorobutadiene	µg	10		6.4
Hexachlorocyclopentadiene	µg	ND	U	ND U
Hexachloroethane	µg	ND	U	ND U
Indeno(1,2,3-c,d)pyrene	µg	ND	U	ND U
Isophorone	µg	2.8		41

Table 3.17
SBPA and Off-Site ISVE System Results
for Method TO-13 (SVOCs) - November 2006
American Chemical Service
Griffith, Indiana

Compounds	Units	Sampled 11/2/2006			
		SBPA ISVE		Off-Site ISVE	
Naphthalene	µg	17		98	
Nitrobenzene	µg	ND	U	ND	U
N-Nitroso-di-n-propylamine	µg	ND	U	ND	U
N-Nitrosodiphenylamine	µg	ND	U	ND	U
Pentachlorophenol	µg	ND	U	ND	U
Phenanthrene	µg	ND	U	ND	U
Phenol	µg	ND	U	ND	U
Pyrene	µg	ND	U	ND	U
Total	µg	91.30		268.20	

Notes:

µg - Microgram

NC - Not calculated

ND - Non-detect

Qualifiers:

J - Result is estimated

U - below reported quantitation limit

/ - Laboratory data qualifier

/ - Data validation qualifier

Table 3.18
SBPA and Off-Site ISVE System Results
for Method TO-13 (SVOCs) - December 2006
American Chemical Service
Griffith, Indiana

Compounds	Units	Sampled 12/11/06		
		SBPA ISVE	Off-Site ISVE	
1,2,4-Trichlorobenzene	µg	19	U	3.6
1,2-Dichlorobenzene	µg	28	U	65
1,3-Dichlorobenzene	µg	3.4	U	2.5
1,4-Dichlorobenzene	µg	6.9	U	7.9
2,4,5-Trichlorophenol	µg	ND	U	ND
2,4,6-Trichlorophenol	µg	ND	U	ND
2,4-Dichlorophenol	µg	ND	U	ND
2,4-Dimethylphenol	µg	ND	U	ND
2,4-Dinitrophenol	µg	ND	U	ND
2,4-Dinitrotoluene	µg	ND	U	ND
2,6-Dinitrotoluene	µg	ND	U	ND
2-Chloronaphthalene	µg	ND	U	ND
2-Chlorophenol	µg	ND	U	ND
2-Methylnaphthalene	µg	16	U	16
2-Methylphenol (o-Cresol)	µg	ND	U	ND
2-Nitroaniline	µg	ND	U	ND
2-Nitrophenol	µg	ND	U	ND
3,3'-Dichlorobenzidine	µg	ND	U	ND
3-Nitroaniline	µg	ND	U	ND
4,6-Dinitro-2-methylphenol	µg	ND	U	ND
4-Bromophenyl-phenyl Ether	µg	ND	U	ND
4-Chloro-3-methylphenol	µg	ND	U	ND
4-Chloroaniline	µg	ND	U	ND
4-Chlorophenyl-phenyl Ether	µg	ND	U	ND
4-Methylphenol/3-Methylphenol	µg	ND	U	3.3
4-Nitroaniline	µg	ND	U	ND
4-Nitrophenol	µg	ND	U	ND
Acenaphthene	µg	ND	U	ND
Acenaphthylene	µg	ND	U	ND
Anthracene	µg	ND	U	ND
Benzo(a)anthracene	µg	ND	U	ND
Benzo(a)pyrene	µg	ND	U	ND
Benzo(b)fluoranthene	µg	ND	U	ND
Benzo(g,h,i)perylene	µg	ND	U	ND
Benzo(k)fluoranthene	µg	ND	U	ND
bis(2-Chloroethoxy) Methane	µg	ND	U	ND
bis(2-Chloroethyl) Ether	µg	3.2	U	6.2
bis(2-Ethylhexyl)phthalate	µg	1.8	J/J	3.5
Butylbenzylphthalate	µg	ND	U	ND
Chrysene	µg	ND	U	ND
Dibenz(a,h)anthracene	µg	ND	U	ND
Dibenzofuran	µg	ND	U	ND
Diethylphthalate	µg	ND	U	ND
Dimethylphthalate	µg	ND	U	ND
di-n-Butylphthalate	µg	ND	U	ND
Di-n-Octylphthalate	µg	ND	U	ND
Fluoranthene	µg	ND	U	ND
Fluorene	µg	ND	U	ND
Hexachlorobenzene	µg	ND	U	ND
Hexachlorobutadiene	µg	13	U	4.8
Hexachlorocyclopentadiene	µg	ND	U	ND
Hexachloroethane	µg	ND	U	ND
Indeno(1,2,3-c,d)pyrene	µg	ND	U	ND
Isophorone	µg	3	U	42

Table 3.18
SBPA and Off-Site ISVE System Results
for Method TO-13 (SVOCs) - December 2006
American Chemical Service
Griffith, Indiana

Compounds	Units	Sampled 12/11/06			
		SBPA ISVE	Off-Site ISVE		
Naphthalene	µg	26	U	76	U
Nitrobenzene	µg	ND	U	ND	U
<i>N</i> -Nitroso-di- <i>n</i> -propylamine	µg	ND	U	ND	U
<i>N</i> -Nitrosodiphenylamine	µg	ND	U	ND	U
Pentachlorophenol	µg	ND	U	ND	U
Phenanthrene	µg	ND	U	ND	U
Phenol	µg	ND	U	5.2	U
Pyrene	µg	ND	U	ND	U
Total	µg	120.30		236.00	

Notes:

µg - Microgram
 NC - Not calculated
 ND - Non-detect

Qualifiers:

J - Result is estimated
 U - below reported quantitation limit
 _/ - Laboratory data qualifier
 /_ - Data validation qualifier

Table 3.19
Off-Site In-Situ Vapor Extraction (ISVE) System Well Monitoring Data
Fourth Quarter 2006
American Chemical Service NPL Site
Griffith, Indiana

Well ID	Date	Flow (cfm)	Vac (" H ₂ O)	VOCs (ppm)	Comments
SVE-01	10/19/2006	Water	84.0	181	
	11/29/2006	Water	78.0	105	
	12/21/2006	Water	88.5	92	
SVE-02	10/19/2006	12	84.0	129	
	11/29/2006	14	79.5	47	
	12/21/2006	9	89.5	82	
SVE-03	10/19/2006	Water	84.5	112	
	11/29/2006	Water	79.0	51	
	12/21/2006	119	88.5	74	
SVE-04	10/19/2006	8	84.0	143	
	11/29/2006	12	79.0	55	
	12/21/2006	11	90.0	75	
SVE-05	10/19/2006	14	84.0	159	
	11/29/2006	13	78.5	67	
	12/21/2006	11	89.5	77	
SVE-06	10/19/2006	Water	84.0	104	
	11/29/2006	Water	79.0	59	
	12/21/2006	Water	89.0	93	
SVE-07	10/19/2006	Water	83.5	119	
	11/29/2006	Water	78.0	52	
	12/21/2006	Water	85.5	Water	
SVE-08	10/19/2006	Water	83.5	Water	
	11/29/2006	Water	73.5	23	
	12/21/2006	4	9.0	52	
SVE-09	10/19/2006	160	83.5	81	
	11/29/2006	197	77.5	18	
	12/21/2006	Water	79.0	53	
SVE-10	10/19/2006	91	83.5	188	
	11/29/2006	Water	78.0	30	
	12/21/2006	130	82.5	93	
SVE-11	10/19/2006	Water	70.0	105	
	11/29/2006	Water	63.5	28	
	12/21/2006	Water	81.5	63	
SVE-12	10/19/2006	15	84.0	130	
	11/29/2006	Water	78.5	43	
	12/21/2006	Water	89.5	71	
SVE-13	10/19/2006	12	82.5	503	
	11/29/2006	13	77.0	636	
	12/21/2006	7	87.5	209	
SVE-14	10/19/2006	Water	83.5	Water	
	11/29/2006	Water	77.5	1344	
	12/21/2006	7	89.0	3090	
SVE-15	10/19/2006	260	83.5	728	
	11/29/2006	Water	56.0	372	Vacuum fluctuated between 51-61
	12/21/2006	Water	88.0	407	

Table 3.19
Off-Site In-Situ Vapor Extraction (ISVE) System Well Monitoring Data
Fourth Quarter 2006
American Chemical Service NPL Site
Griffith, Indiana

Well ID	Date	Flow (cfm)	Vac ($\text{" H}_2\text{O}$)	VOCs (ppm)	Comments
SVE-16	10/19/2006	Water	83.0	1705	
	11/29/2006	8	45.5	1021	
	12/21/2006	Water	87.5	1386	
SVE-17	10/19/2006	15	83.5	199	
	11/29/2006	Water	78.5	97	
	12/21/2006	10	89.0	143	
SVE-18	10/19/2006	9	83.5	624	
	11/29/2006	18	77.5	105	
	12/21/2006	5	87.5	1713	
SVE-19	10/19/2006	Water	83.5	106	
	11/29/2006	Water	79.0	76	
	12/21/2006	Water	88.0	115	
SVE-20	10/19/2006	Water	83.5	95	
	11/29/2006	Water	77.0	64	
	12/21/2006	Water	87.5	115	
SVE-21	10/19/2006	115	83.5	141	
	11/29/2006	36	63.0	83	Vacuum fluctuated between 58-68
	12/21/2006	30	82.5	179	
SVE-22	10/19/2006	20	76.5	4448	
	11/29/2006	30	68.0	1190	
	12/21/2006	Water	87.5	1377	
SVE-23	10/19/2006	Water	79.0	1671	
	11/29/2006	Water	69.5	1256	
	12/21/2006	Water	88.0	1482	
SVE-24	10/19/2006	Water	72.0	1506	
	11/29/2006	160	61.0	1056	
	12/21/2006	178	87.5	1454	
SVE-25	10/19/2006	150	76.5	1271	
	11/29/2006	Water	68.0	728	
	12/21/2006	Water	86.5	697	
SVE-26	10/19/2006	37	83.5	366	
	11/29/2006	Water	78.0	153	
	12/21/2006	Water	87.5	288	
SVE-27	10/19/2006	10	52.5	276	
	11/29/2006	9	43.0	102	
	12/21/2006	3	50.0	154	
SVE-28	10/19/2006	20	83.5	349	
	11/29/2006	19	77.5	299	
	12/21/2006	4	87.5	135	
SVE-29	10/19/2006	45	83.5	552	
	11/29/2006	Water	78.0	315	
	12/21/2006	Water	87.5	Water	
SVE-30	10/19/2006	12	83.5	2303	
	11/29/2006	10	78.0	1348	
	12/21/2006	4	88.0	1448	

Table 3.19
Off-Site In-Situ Vapor Extraction (ISVE) System Well Monitoring Data
Fourth Quarter 2006
American Chemical Service NPL Site
Griffith, Indiana

Well ID	Date	Flow (cfm)	Vac ($\text{"H}_2\text{O}$)	VOCs (ppm)	Comments
SVE-31	10/19/2006	13	82.5	980	
	11/29/2006	17	78.0	37	
	12/21/2006	7	87.5	23	
SVE-32	10/19/2006	Water	81.0	Water	
	11/29/2006	12	28.0	255	
	12/21/2006	15	42.5	21	
SVE-33	10/19/2006	16	83.5	155	
	11/29/2006	13	79.0	459	
	12/21/2006	7	87.5	44	
SVE-34	10/19/2006	78	83.5	852	
	11/29/2006	Water	78.5	463	
	12/21/2006	100	87.0	418	
SVE-35	10/19/2006	18	83.0	86	
	11/29/2006	15	77.5	51	
	12/21/2006	6	87.5	41	
SVE-36	10/19/2006	9	82.5	368	
	11/29/2006	13	77.0	2338	
	12/21/2006	8	87.5	100	
SVE-37	10/19/2006	Water	82.0	117	
	11/29/2006	Water	75.0	60	
	12/21/2006	7	87.0	25	
SVE-38	10/19/2006	Water	83.0	1411	
	11/29/2006	Water	77.0	453	
	12/21/2006	Water	87.5	354	
SVE-39	10/19/2006	15	83.0	102	
	11/29/2006	16	78.5	52	
	12/21/2006	5	87.0	33	
SVE-40	10/19/2006	Water	56.5	1217	
	11/29/2006	Water	45.5	388	
	12/21/2006	Water	87.0	304	
SVE-41	10/19/2006	Water	83.0	955	
	11/29/2006	Water	77.5	473	
	12/21/2006	Water	87.5	415	
SVE-42	10/19/2006	10	82.5	675	
	11/29/2006	21	77.5	73	
	12/21/2006	11	86.5	123	

Notes:

"Water" - water present in vapor stream, preventing data collection

Beginning in March 2006, flow is measured using a VelociCalc 8384 flow meter.

Differential pressure is no longer measured.

Beginning in August 2006, vacuum pressures are measured with an Extech Manometer Model 407910.

Table 3.20
Off-Site In-Situ Vapor Extraction (ISVE) System Header Monitoring Data - Fourth Quarter 2006
American Chemical Service NPL Site
Griffith, Indiana

Date	KP1 Line Press (psia)	KP1 Flow (scfm)	KP1 Vac (" H ₂ O)	KP2 Line Press (psia)	KP2 Flow (scfm)	KP2 Vac (" H ₂ O)	OFCA1 Vac (" H ₂ O)	OFCA2 Vac (" H ₂ O)	OFCA3 Vac (" H ₂ O)	Dilution Flow (cfm)	Blower Inf Line Press (psia)	Blower Inf Flow (scfm)
10/19/2006	11.6	0	84	11.7	0	83.5	83.5	83	84	0	11.5	0
11/29/2006	11.8	349	79.5	11.8	0	79.5	79	78	79	0	11.7	694
12/21/2006	11.5	0	90	11.5	0	90	88.5	88.5	89.5	0	11.4	687

Date	Blower Inf Vac (" H ₂ O)	Blower Inf VOC (ppm)	Blower Inf Temp. (°F)	Blower Eff Line Press (psia)	Blower Eff Flow (scfm)	Blower Eff Press (" H ₂ O)	Blower Eff VOC (ppm)	Blower Eff Temp. (°F)	Filter Diff Press (" H ₂ O)	Ambient Temperature (°F)	Barometric Pressure ("Hg)	Humidity (%)
10/19/2006	89	459	68	15.5	525	24.0	-	152	7.0	45	29.87	66%
11/29/2006	83	61	66	15.7	647	28.5	-	150	7.0	60	29.91	88%
12/21/2006	94	330	62	15.7	527	-26.5	-	154	6.0	46	30.06	93%

Notes:

"-" = data not collected

cfm = cubic feet per minute

" H₂O = inches of water

ppm = parts per million

VOCs = volatile organic compounds

psia = pounds per square inch, atmosphere

" Hg = inches of mercury

°F = degrees Fahrenheit

Table 3.21
SBPA In-Situ Vapor Extraction (ISVE) System Well Monitoring Data
Fourth Quarter 2006
American Chemical Service NPL Site
Griffith, Indiana

Well ID	Date	Flow (cfm)	Vac (in H ₂ O)	VOCs (ppm)	Comments
SVE-43	10/19/2006	22	81.0	28	
	11/29/2006	22	94.5	71	
	12/21/2006	0.1	1.0	44	
SVE-44	10/19/2006	80	79.5	21	
	11/29/2006	Water	93.5	6	
	12/21/2006	45	18.5	78	
SVE-45	10/19/2006	14	80.5	38	
	11/29/2006	32	94.0	127	
	12/21/2006	22	115.0	49	
SVE-46	10/19/2006	60	81.0	64	
	11/29/2006	25	94.0	59	
	12/21/2006	Water	114.5	60	
SVE-47	10/19/2006	34	81.0	642	
	11/29/2006	19	94.0	635	
	12/21/2006	23	114.0	194	
SVE-48	10/19/2006	Water	79.5	650	
	11/29/2006	Water	93.0	523	
	12/21/2006	Water	87.5	142	
SVE-49	10/19/2006	20	81.5	60	
	11/29/2006	28	94.5	183	
	12/21/2006	15	-	-	Air injection well
SVE-50	10/19/2006	14	-	-	Air injection well
	11/29/2006	16	-	-	Air injection well
	12/21/2006	Water	115.0	104	
SVE-51	10/19/2006	Water	79.5	81	
	11/29/2006	20	94.0	17	
	12/21/2006	33	-	-	Air injection well
SVE-52	10/19/2006	-	-	-	
	11/29/2006	-	-	-	
	12/21/2006	-	-	-	
SVE-53	10/19/2006	-	-	-	
	11/29/2006	-	-	-	
	12/21/2006	-	-	-	
SVE-54	10/19/2006	15	-	-	Air injection well
	11/29/2006	17	-	-	Air injection well
	12/21/2006	Water	106.0	185	
SVE-55	10/19/2006	17	81.5	427	
	11/29/2006	28	94.5	288	
	12/21/2006	46	115.5	227	
SVE-56	10/19/2006	13	81.5	197	
	11/29/2006	25	95.5	99	
	12/21/2006	54	114.5	84	
SVE-57	10/19/2006	30	70.0	232	
	11/29/2006	22	94.0	117	
	12/21/2006	23	111.5	74	
SVE-58	10/19/2006	32	75.0	332	
	11/29/2006	Water	98.0	260	
	12/21/2006	Water	114.5	118	

Table 3.21
SBPA In-Situ Vapor Extraction (ISVE) System Well Monitoring Data
Fourth Quarter 2006
American Chemical Service NPL Site
Griffith, Indiana

Well ID	Date	Flow (cfm)	Vac (in H ₂ O)	VOCs (ppm)	Comments
SVE-59	10/19/2006	Water	69.0	5	
	11/29/2006	Water	93.5	209	
	12/21/2006	Water	114.5	85	
SVE-60	10/19/2006	13	83.5	234	
	11/29/2006	30	98.0	258	
	12/21/2006	Water	114.0	101	
SVE-61	10/19/2006	-	-	-	
	11/29/2006	-	-	-	
	12/21/2006	-	-	-	
SVE-62	10/19/2006	-	-	-	
	11/29/2006	-	-	-	
	12/21/2006	-	-	-	
SVE-63	10/19/2006	12	81.0	686	
	11/29/2006	Water	95.0	310	
	12/21/2006	33	116.0	172	
SVE-64	10/19/2006	18	80.0	1498	
	11/29/2006	26	95.0	651	
	12/21/2006	66	115.0	192	
SVE-65	10/19/2006	18	79.0	2564	
	11/29/2006	Water	94.0	757	
	12/21/2006	22	-	-	Air injection well
SVE-66	10/19/2006	24	79.5	228	
	11/29/2006	20	95.0	147	
	12/21/2006	30	114.0	73	
SVE-67	10/19/2006	38	61.0	Water	
	11/29/2006	45	62.5	530	
	12/21/2006	Water	113.0	55	
SVE-68	10/19/2006	25	80.5	227	
	11/29/2006	Water	95.5	147	
	12/21/2006	Water	76.0	45	
SVE-69	10/19/2006	Water	55.0	Water	
	11/29/2006	21	29.0	17	
	12/21/2006	34	93.5	56	
SVE-70	10/19/2006	84	84.0	49	
	11/29/2006	NM	98.0	25	Blockage - could not measure flow
	12/21/2006	NM	22.5	61	Blockage - could not measure flow
SVE-71	10/19/2006	16	83.5	410	
	11/29/2006	Water	98.5	539	
	12/21/2006	46	-	-	Air injection well
SVE-72	10/19/2006	-	-	-	
	11/29/2006	-	-	-	
	12/21/2006	-	-	-	
SVE-73	10/19/2006	17	-	-	Air injection well
	11/29/2006	23	-	-	Air injection well
	12/21/2006	29	106.0	99	
SVE-74	10/19/2006	15	83.5	2824	
	11/29/2006	41	98.0	1983	
	12/21/2006	55	121.5	144	

Table 3.21
SBPA In-Situ Vapor Extraction (ISVE) System Well Monitoring Data
Fourth Quarter 2006
American Chemical Service NPL Site
Griffith, Indiana

Well ID	Date	Flow (cfm)	Vac ($\text{" H}_2\text{O}$)	VOCs (ppm)	Comments
SVE-75	10/19/2006	Water	82.5	159	
	11/29/2006	Water	95.0	104	
	12/21/2006	Water	95.0	54	
SVE-76	10/19/2006	28	81.0	100	
	11/29/2006	33	95.0	60	
	12/21/2006	71	110.0	61	
SVE-77	10/19/2006	-	-	-	
	11/29/2006	-	-	-	
	12/21/2006	-	-	-	
SVE-78	10/19/2006	-	-	-	
	11/29/2006	-	-	-	
	12/21/2006	-	-	-	
SVE-79	10/19/2006	19	-	-	Air injection well
	11/29/2006	27	-	-	Air injection well
	12/21/2006	130	119.0	85	
SVE-80	10/19/2006	22	84.0	108	
	11/29/2006	Water	98.0	301	
	12/21/2006	Water	117.5	89	
SVE-81	10/19/2006	29	-	-	Air injection well
	11/29/2006	28	-	-	Air injection well
	12/21/2006	Water	118.0	195	
SVE-82	10/19/2006	Water	83.5	307	
	11/29/2006	Water	98.0	317	
	12/21/2006	24	-	-	Air injection well
SVE-83	10/19/2006	14	85.0	1871	
	11/29/2006	Water	98.0	376	
	12/21/2006	Water	112.0	176	
SVE-84	10/19/2006	10	84.0	157	
	11/29/2006	Water	98.5	720	
	12/21/2006	-	-	-	
SVE-85	10/19/2006	13	84.0	1700	
	11/29/2006	25	98.0	2421	
	12/21/2006	Water	120.0	229	
SVE-86	10/19/2006	14	84.0	531	
	11/29/2006	17	98.5	1246	
	12/21/2006	43	116.0	110	
SVE-87	10/19/2006	13	59.0	791	
	11/29/2006	-	-	-	
	12/21/2006	Water	109.0	144	
SVE-88	10/19/2006	-	-	-	
	11/29/2006	-	-	-	
	12/21/2006	-	-	-	

Notes:

"-" = data not collected

"Water" - water present in vapor stream, preventing data collection

NM = Not measured, reason given in comments column

Beginning in March 2006, flow is measured using a VelociCalc 8384 flow meter.

Differential pressure is no longer measured.

Beginning in August 2006, vacuum pressures are measured with an Extech Manometer

Model 407910.

Table 3.22
SBPA In-Situ Vapor Extraction (ISVE) System Header Monitoring Data - Fourth Quarter 2006
American Chemical Service NPL Site
Griffith, Indiana

Date	Line Press (psia)	Flow	Vac	Line Press (psia)	Flow	Vac	Dilution	Blower Inf Line Press (psia)	Blower Inf Flow (scfm)	Blower Inf Vac (" H ₂ O)	Blower Inf VOC (ppm)
		(scfm)	(" H ₂ O)		(scfm)	(" H ₂ O)	(cfm)				
10/19/2006	11.6	1559	85	11.6	0	86	0	11.1	1522	100	-
11/29/2006	11.2	1326	97	11.2	0	98.5	0	11.1	1669	100	325
12/21/2006	10.6	0	116	10.5	1444	120.0	0	11.2	1678	100	-

Date	Blower Inf Temp. (°F)	Blower Eff Line Press (psia)	Blower Eff Flow (scfm)	Blower Eff Press (" H ₂ O)	Blower Eff VOC (ppm)	Blower Eff Temp. (°F)	Filter Diff Press (" H ₂ O)	Ambient Temperature (°F)	Barometric Pressure ("Hg)	Humidity (%)
10/19/2006	59	14.8	892	4.0	345	147	11.0	43	29.93	70%
11/29/2006	58	14.7	781	0.0	375	194	20.0	62	29.94	82%
12/21/2006	58	14.8	936	0.0	-	182	20.0	43	30.18	87%

Notes:

"—" = data not collected

scfm = cubic feet per minute

" H₂O = inches of water

ppm = parts per million

VOCs = volatile organic compounds

psia = pounds per square inch, atmosphere

" Hg = inches of mercury

°F = degrees Fahrenheit

Table 3.23
Schedule of Product Removal Activities - Fourth Quarter 2006
American Chemical Service
Griffith, Indiana

Date	Well	Amount of Product Removed
October 24, 2006	SVE-53	10 gallons
October 30, 2006	SVE-72	8 gallons
November 6, 2006	SVE-53	10 gallons
November 14, 2006	SVE-72	4 gallons
November 27, 2006	SVE-53	6 gallons
Total Product Removed		38 gallons

Table 6.1
Water Table Elevations Across the Barrier Wall and Near the PGCS - Fourth Quarter 2006
American Chemical Service NPL Site
Griffith, Indiana

Upper Aquifer Wells

Well Designation	Reference Points			12/22/2006		Notes	Difference Across Barrier Wall (if applicable) ¹
	East	North	TOIC	Level	Elevation		
MW11	6377	7329	640.47	4.14	636.33		n/a
MW13	5050	7814	634.08	2.34	631.74		n/a
MW37	5395	7976	636.78	3.37	633.41		n/a
MW46	4526	7424	633.32	2.28	631.04	Flooded	n/a
MW48	5669	7814	636.36	3.74	632.62		n/a
MW49	5551	7650	637.00	3.37	633.63		n/a

Staff Gauges & Piezometers

Well Designation	Reference Points			12/22/2006		Notes	Difference Across Barrier Wall (if applicable) ¹
	East	North	TOSG	Level	Elevation		
P23	4689	7018	636.18	4.45	631.73		n/a
P25	5131	7510	633.33	1.26	632.07		n/a
P26	4764	7309	634.23	3.08	631.15		n/a
P27	4904	7020	639.70	8.26	631.44		n/a
P28	5883	7486	644.53	9.55	634.98		n/a
P32	5746	7026	642.32	10.49	631.83		n/a
P40	5931	7241	638.77	3.08	635.69		n/a
P41	5663	7377	637.23	1.65	635.58		n/a
P49	5145	6949	638.98	9.52	629.46		n/a
SG13	4819	7209	631.53	5.38	630.91	TOSG = 6.0' mark	n/a

PGCS Piezometer Sets

Well Designation	Reference Points			12/22/2006		Notes	Difference Across Barrier Wall (if applicable) ¹
	East	North	TOC	Level	Elevation		
P81	5577	7581	636.19	2.71	633.48		n/a
P82	5577	7572	635.77	2.56	633.21		n/a
P83	5577	7561.6	635.95	2.24	633.71		n/a
P84	5322	7603	634.35	2.05	632.30		n/a
P85	5326	7594	634.08	1.63	632.45		n/a
P86	5329	7585	634.41	1.82	632.59		n/a
P87	5121	7466	633.88	3.64	630.24		n/a
P88	5130	7460	633.90	2.01	631.89		n/a
P89	5137	7454	634.02	2.18	631.84		n/a
P90	4881	7152	634.45	3.13	631.32		n/a
P91	4889	7145	634.59	3.82	630.77		n/a
P92	4896	7138.1	633.87	2.96	630.91		n/a

Table 6.1
Water Table Elevations Across the Barrier Wall and Near the PGCS - Fourth Quarter 2006
American Chemical Service NPL Site
Griffith, Indiana

BWES Water Level and Piezometer Pairs

Well Designation	Reference Points			12/22/2006		Notes	Difference Across Barrier Wall (if applicable) ¹
	East	North	TOC	Level	Elevation		
P93R - Outside BW	TBD	TBD	639.05	6.45	632.60	Installed Nov. 2004	-3.09
P94R - Inside BW	TBD	TBD	640.99	11.48	629.51	Installed Nov. 2004	
P95 - Outside BW	5146	6532	638.58	4.04	634.54		-10.38
P96 - Inside BW	5156	6537	641.26	17.10	624.16		
P105 - Outside BW	5885	6678	638.86	2.15	636.71		
P106 - Inside BW	5871	6685	638.10	8.69	629.41		-7.30
P107 - Outside BW	5766	7339	637.42	2.23	635.19		
P108 - Inside BW	5757	7324	638.13	6.38	631.75		-3.44
P109 - Outside BW	5740	6387	644.30	8.29	636.01		
P110 - Inside BW	5705	6382	647.68	21.27	626.41		-9.60
P111 - Outside BW	5551	5950	650.03	14.81	635.22		
P112 - Inside BW	5525	5960	653.36	28.86	624.50		-10.72
P113 - Inside BW	5309	5693	657.53	32.97	624.56		
ORCPZ102 - Outside BW	5331	5612	652.47	17.92	634.55		-9.99
P114 - Inside BW	5035	5729	653.69	29.46	624.23		
P115 - Outside BW	4970	5708	652.50	17.68	634.82		
P116 - Inside BW	5031	6087	646.26	21.94	624.32		
P117 - Outside BW	5014	6087	643.93	8.13	635.80		-11.48
P118 - Inside BW	5402	6539	645.52	19.11	626.41		n/a

Notes:

All depth measurements and elevations are in units of feet.

Elevation is in feet above mean sea level.

TOIC = top of inner casing

TOC = top of casing

TOSG = top of staff gauge

NM = could not measure (reason given under "Notes" column)

n/a = not applicable

¹ A positive value indicates that the water level is higher inside the barrier wall. A negative value indicates that the water level is lower inside the barrier wall.

Table 6.2
Water Levels Inside Barrier Wall - Fourth Quarter 2006
American Chemical Service NPL Site
Griffith, Indiana

Date	On-Site Area					
	Target Level	P-29	P-31	P-32	P-36	P-49
10/12/2006	629.0	630.4	630.9	629.7	624.9	627.7
10/27/2006	629.0	630.4	630.9	629.7	624.9	627.7
11/10/2006	629.0	630.4	630.9	629.7	624.9	627.7
11/22/2006	629.0	630.4	630.9	631.3	626.1	629.2
12/8/2006	629.0	630.4	630.9	631.8	624.9	629.2

Date	Off-Site Area										
	Target Level	P-96	P-110	P-112	P-113	P-114	P-116	P-118	AS-7	AS-8	AS-9
10/12/2006	626.0	620.5	628.0	626.3	626.0	626.3	625.8	627.1	NM	NM	NM
10/19/2006	626.0	NM	627.99	621.90	632.42						
10/27/2006	626.0	620.5	627.5	625.4	625.7	625.8	625.5	626.7	NM	NM	NM
11/10/2006	626.0	620.5	627.6	625.9	626.1	626.4	626.2	626.7	NM	NM	NM
11/14/2006	626.0	NM	628.47	628.16	627.92						
11/22/2006	626.0	621.0	627.8	626.3	627.5	628.1	628.3	626.7	NM	NM	NM
12/8/2006	626.0	620.5	627.7	625.8	625.9	626.2	625.6	626.7	NM	NM	NM
12/21/2006	626.0	NM	630.14	626.96	627.06						

Notes:

All water level elevations are in feet AMSL.

FIGURES

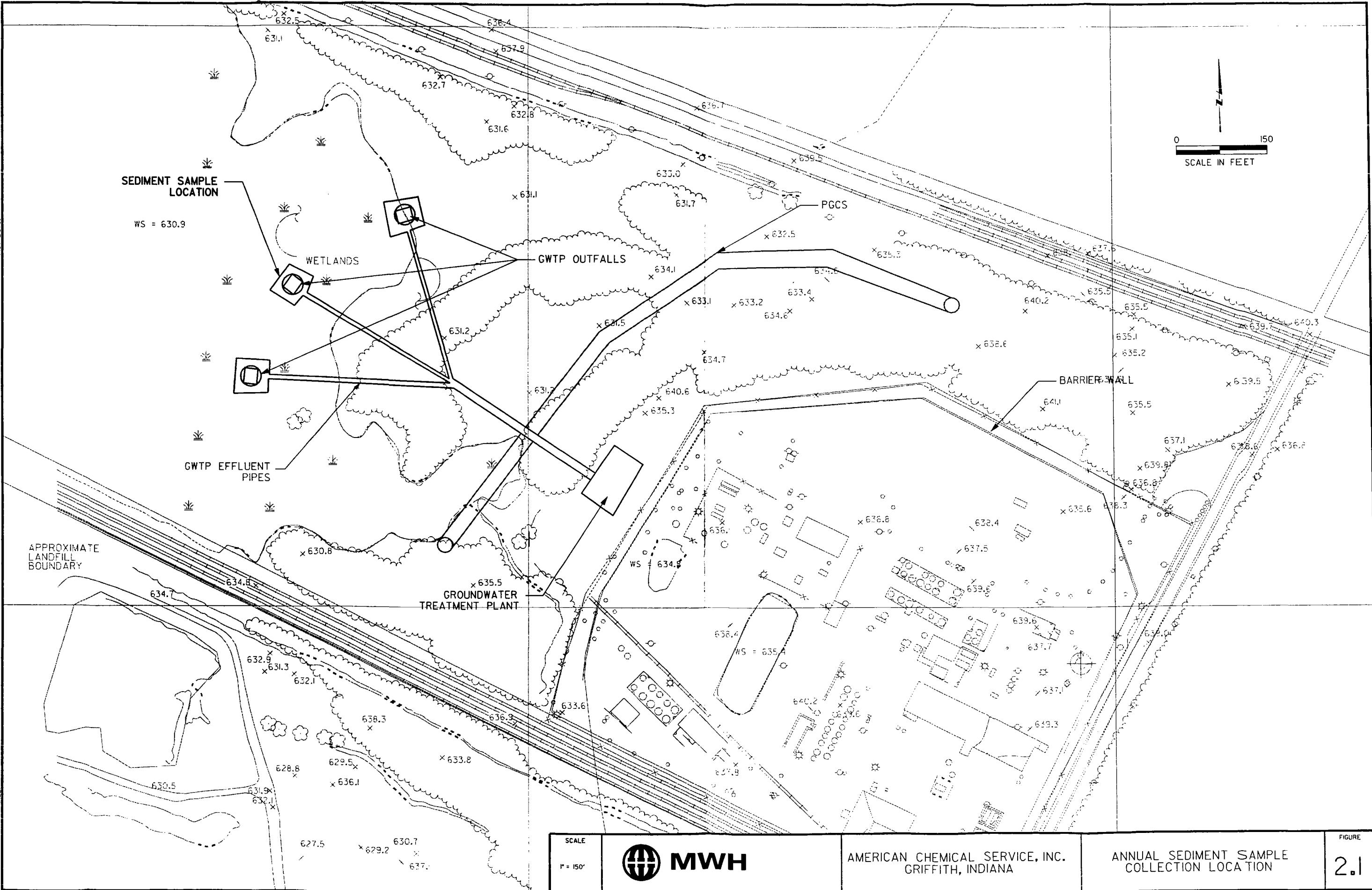


Figure 3.1
VOC Removal Rate
American Chemical Services NPL Site, Griffith, IN

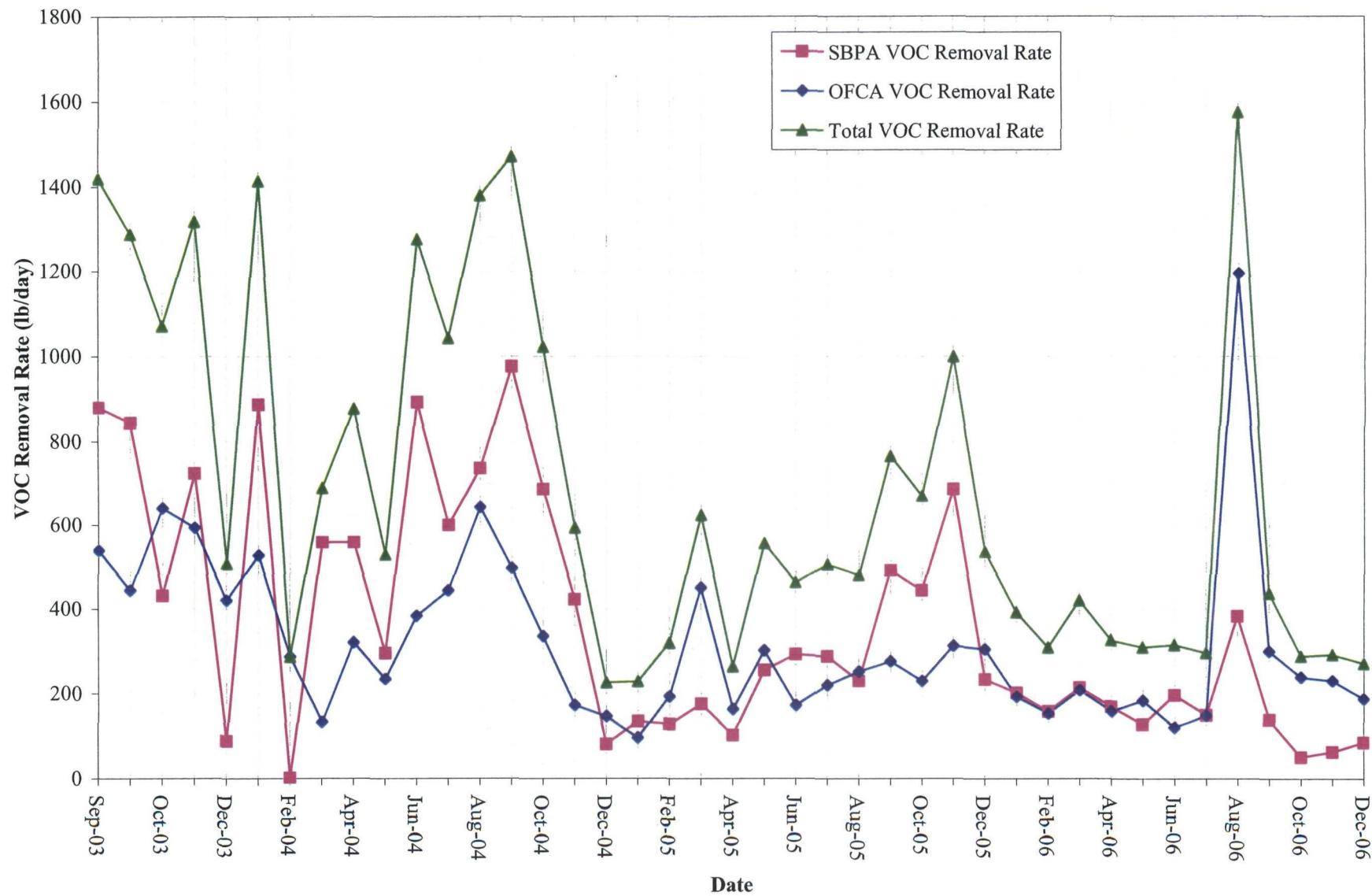
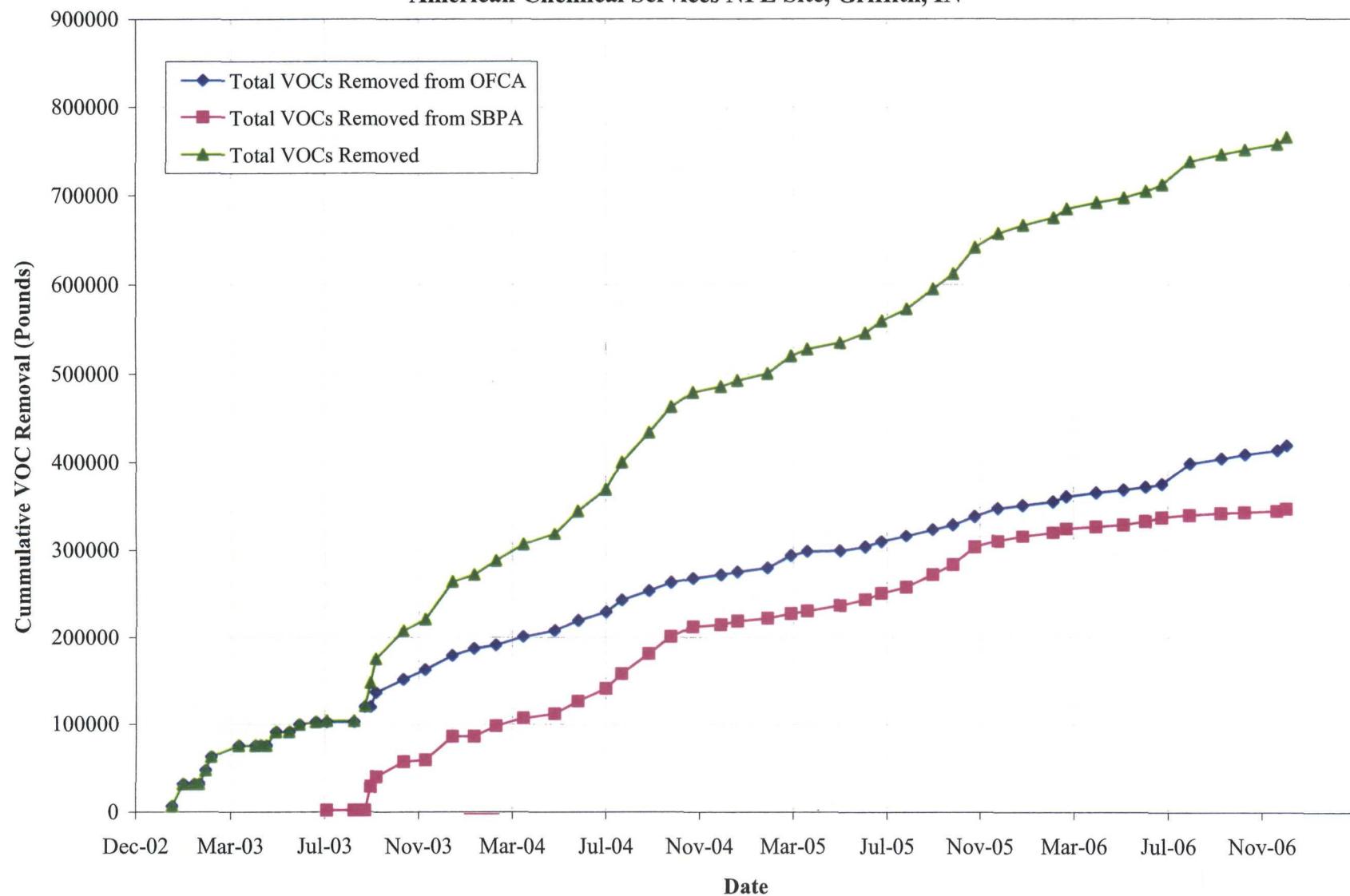
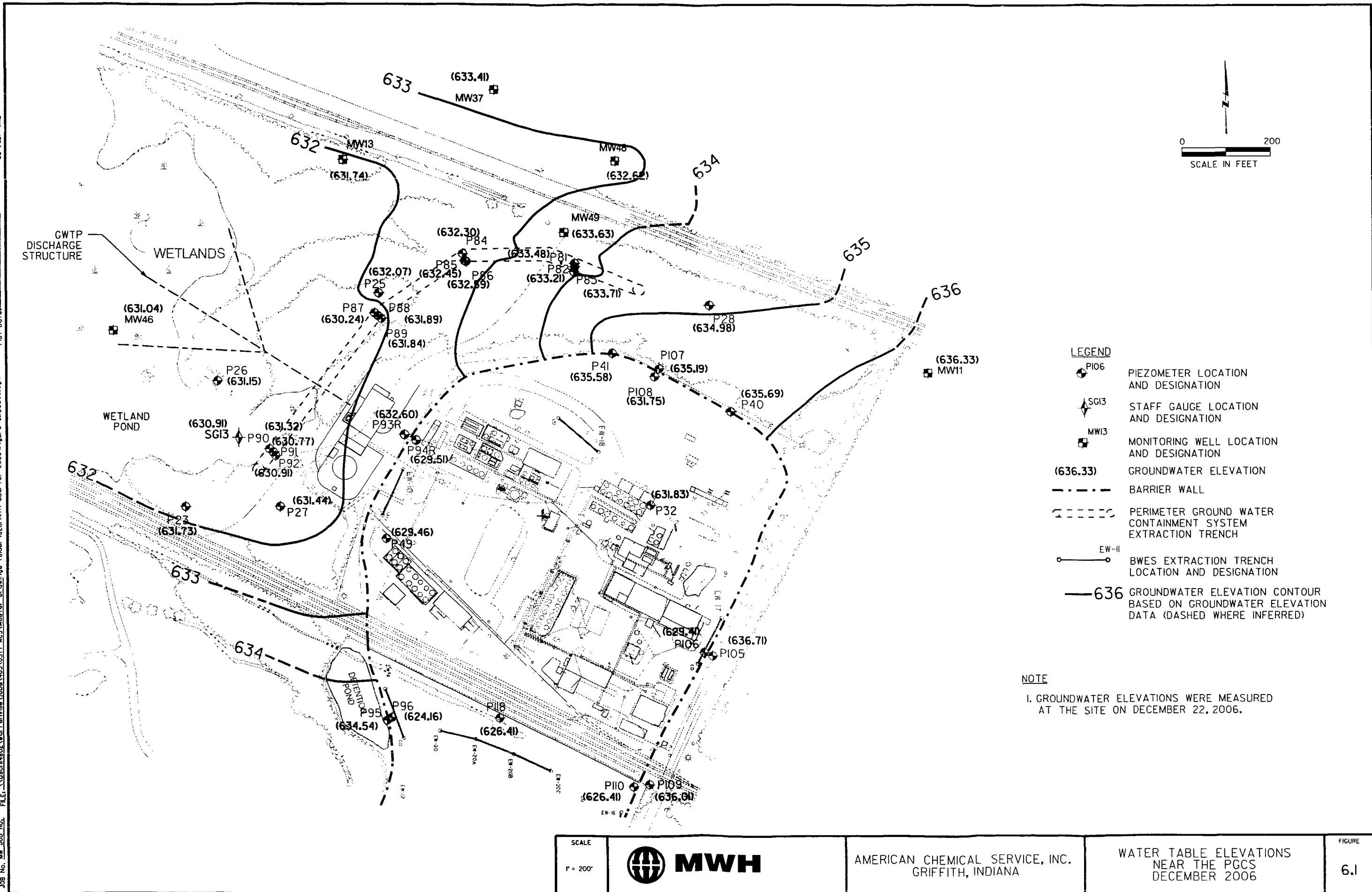
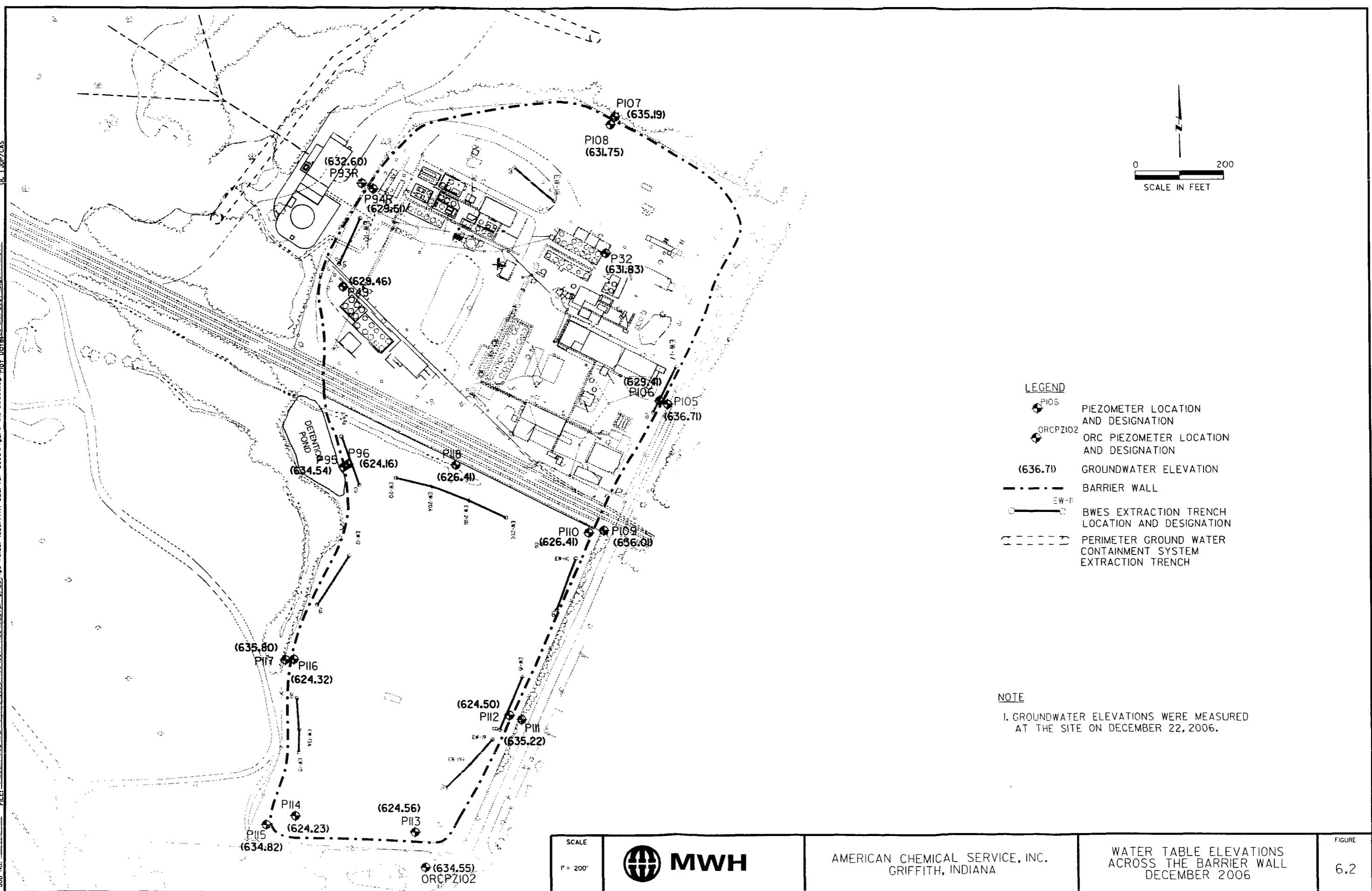
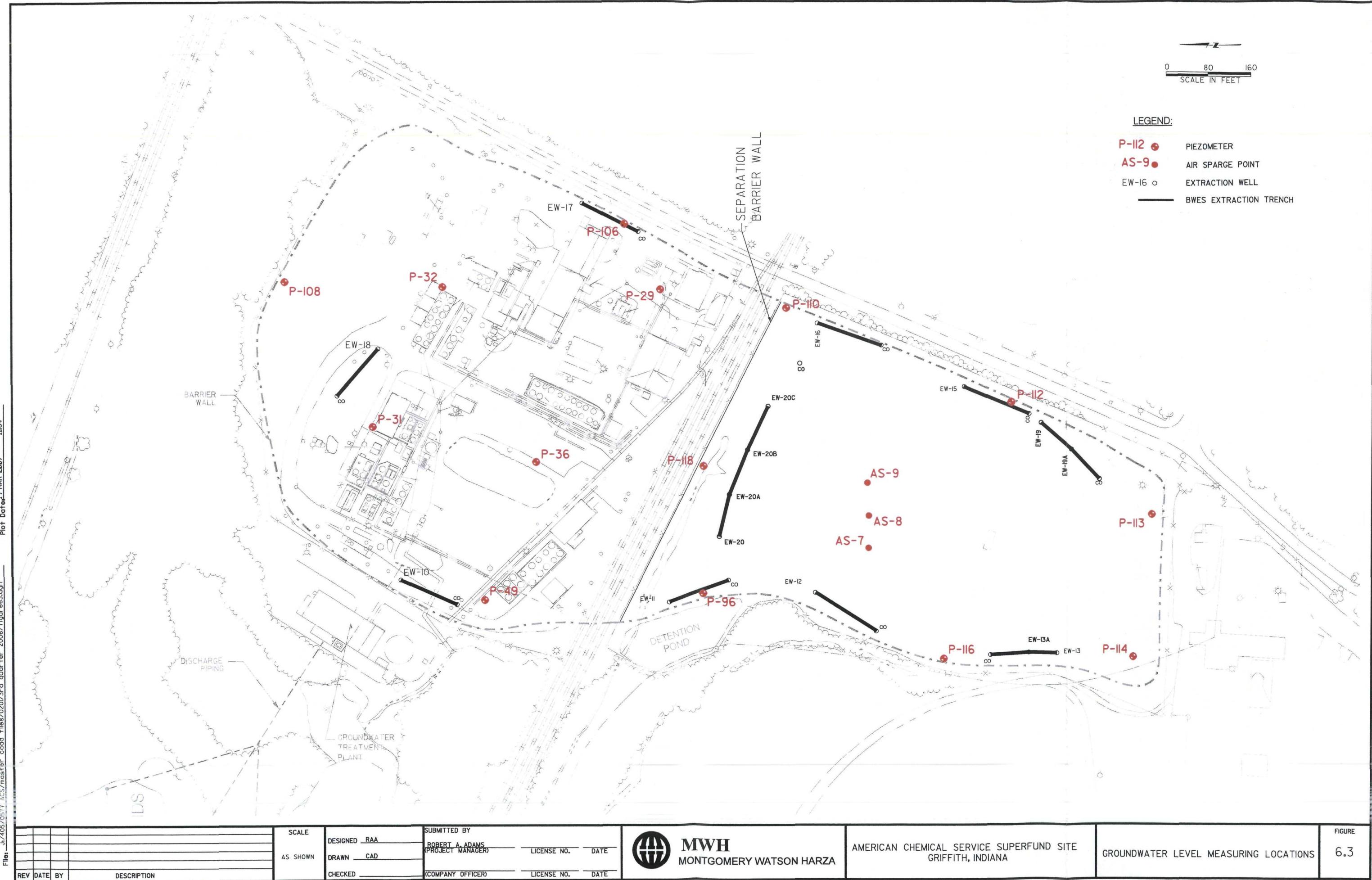


Figure 3.2
Total VOCs Removed
American Chemical Services NPL Site, Griffith, IN









REV	DATE	BY	DESCRIPTION	SCALE	AS SHOWN	SUBMITTED BY	DESIGNED RAA	DRAWN CAD	CHECKED	LICENSE NO.	DATE
						ROBERT A. ADAMS (PROJECT MANAGER)					



MWH
MONTGOMERY WATSON HARZA

AMERICAN CHEMICAL SERVICE SUPERFUND SITE
GRIFFITH, INDIANA

GROUNDWATER LEVEL MEASURING LOCATIONS

FIGURE
6.3

Figure 6.4
Water Level Trends Inside the Barrier Wall (Still Bottoms Pond Area)
ACS NPL Site
Griffith, Indiana

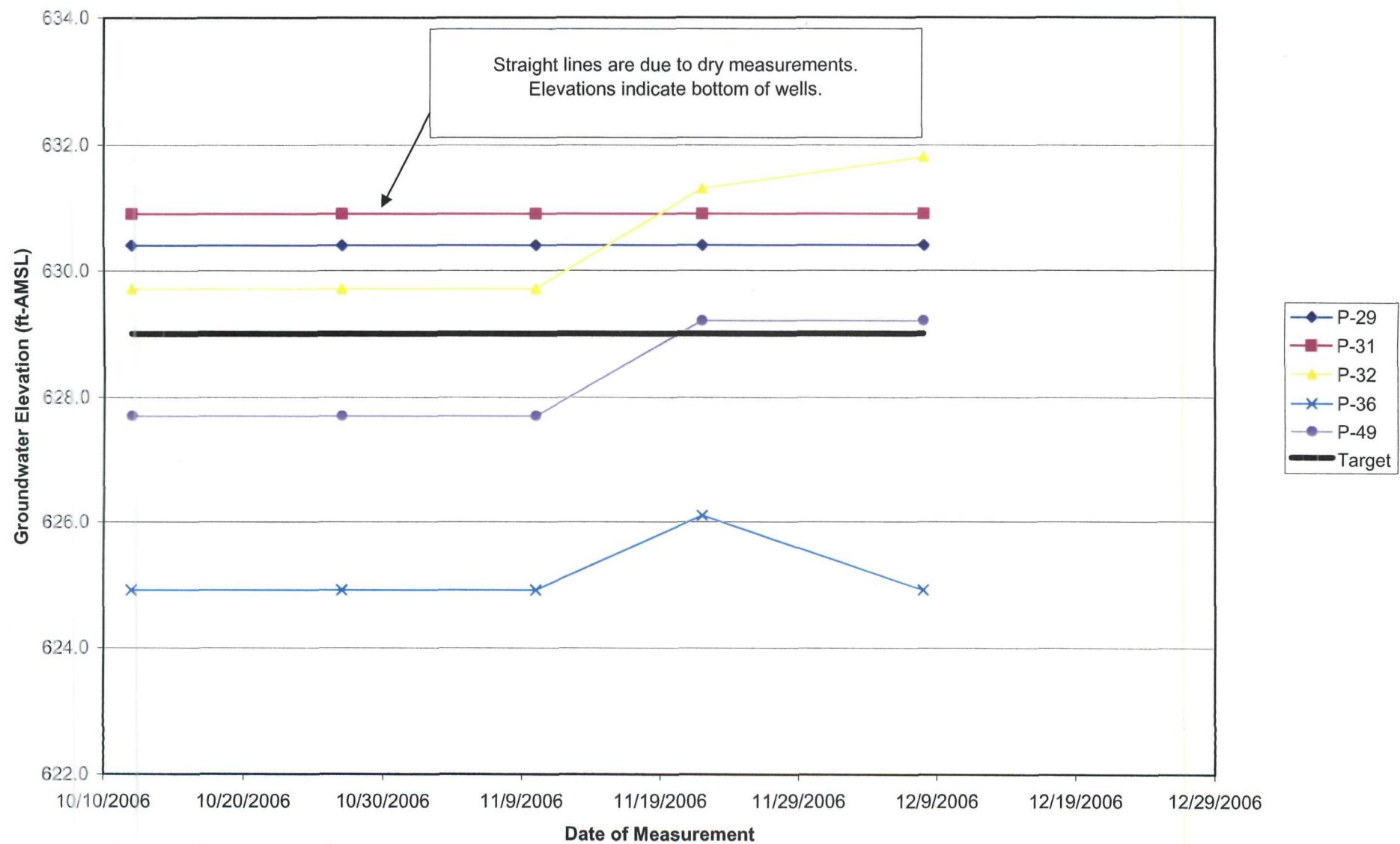
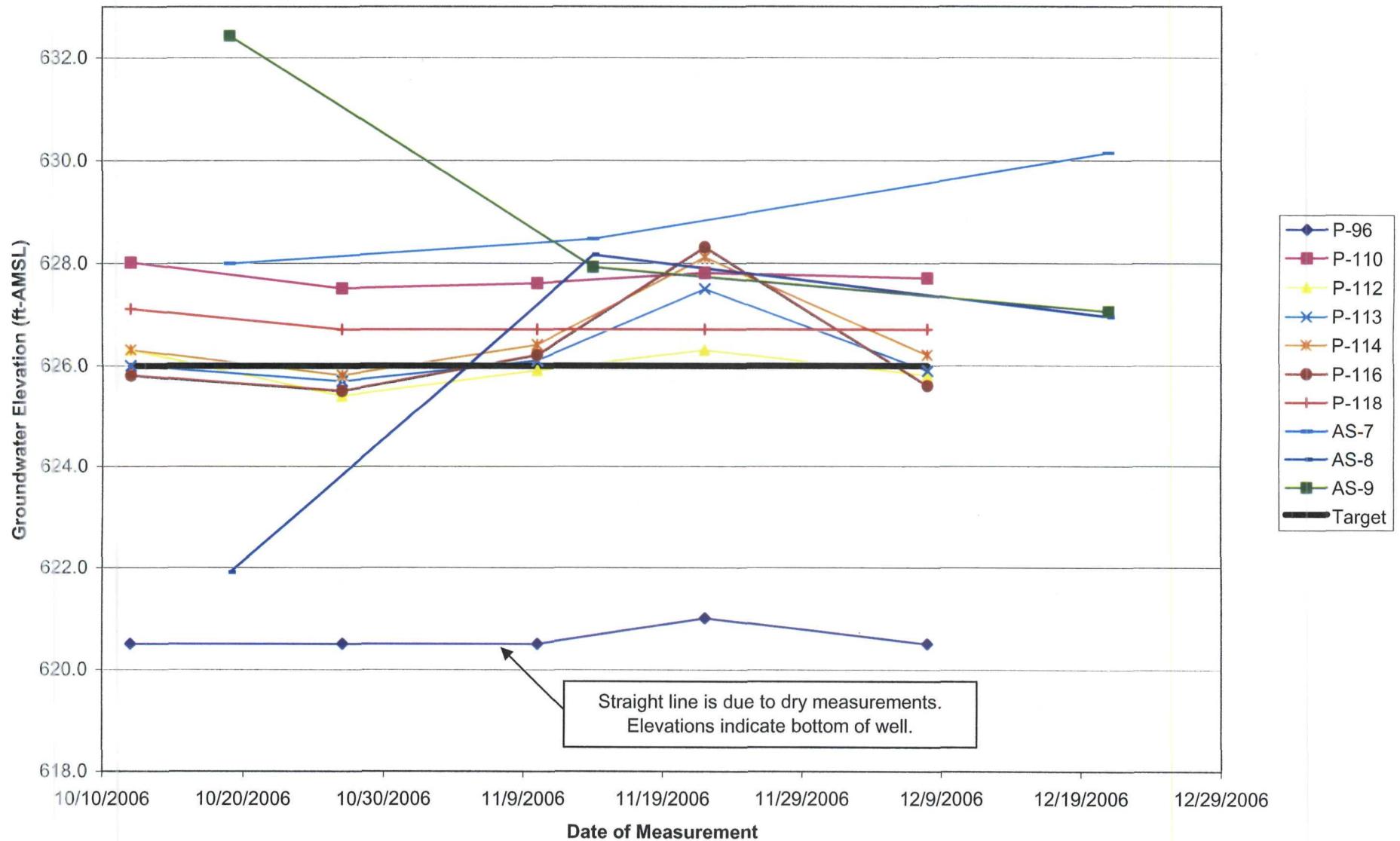


Figure 6.5
Water Level Trends Inside the Barrier Wall (Off-Site Area)
ACS NPL Site
Griffith, Indiana



APPENDIX A

EFFLUENT ANALYTICAL DATA

**October 11, 2006 Compliance Sample
Laboratory Results**

FORM 1
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

Lab Name: COMPUCHEM	Method: 8260B	EFFLUENT
Lab Code: LIBRTY	Case No.:	SAS No.: SDG No.: 11222
Matrix: (soil/water) WATER		Lab Sample ID: 1122201
Sample wt/vol: 25	(g/ml) ML	Lab File ID: 1122201RA61
Level: (low/med)	LOW	Date Received: 10/12/06
% Moisture: not dec.		Date Analyzed: 10/17/06
GC Column: RTX-VMS	ID: 0.18 (mm)	Dilution Factor: 1.0
Soil Extract Volume: _____ (uL)		Soil Aliquot Volume: _____ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) UG/L	Q
74-87-3-----	Chloromethane	0.50	U <i>UJ</i>
75-01-4-----	Vinyl Chloride	0.50	U <i>J</i>
74-83-9-----	Bromomethane	0.50	U <i>UJ</i>
75-00-3-----	Chloroethane	0.48	J <i>J</i>
75-35-4-----	1,1-Dichloroethene	0.50	U <i>UJ</i>
75-15-0-----	Carbon disulfide	0.50	U <i>J</i>
67-64-1-----	Acetone	2.5	U <i>UJ</i>
75-09-2-----	Methylene Chloride	0.45	J <i>J</i>
156-60-5-----	trans-1,2-Dichloroethene	0.50	U <i>UJ</i>
75-34-3-----	1,1-Dichloroethane	0.50	U <i>UJ</i>
156-59-2-----	cis-1,2-Dichloroethene	0.58	J
78-93-3-----	2-butanone	2.5	U <i>UJ</i>
67-66-3-----	Chloroform	0.10	J <i>J</i>
71-55-6-----	1,1,1-Trichloroethane	0.50	U <i>UJ</i>
56-23-5-----	Carbon Tetrachloride	0.50	U <i>J</i>
71-43-2-----	Benzene	0.50	U
107-06-2-----	1,2-Dichloroethane	0.50	U <i>J</i>
79-01-6-----	Trichloroethene	0.50	U <i>UJ</i>
78-87-5-----	1,2-Dichloropropane	0.50	U
75-27-4-----	Bromodichloromethane	0.50	U
10061-01-5-----	cis-1,3-Dichloropropene	0.50	U
108-10-1-----	4-Methyl-2-pentanone	2.5	U
108-88-3-----	Toluene	0.50	U <i>UJ</i>
10061-02-6-----	trans-1,3-Dichloropropene	0.50	U <i>UJ</i>
79-00-5-----	1,1,2-Trichloroethane	0.50	U
127-18-4-----	Tetrachloroethene	0.50	U
591-78-6-----	2-hexanone	2.5	U
124-48-1-----	Dibromochloromethane	0.50	U
108-90-7-----	Chlorobenzene	0.50	U
100-41-4-----	Ethylbenzene	0.50	U
108-38-3-----	m,p-Xylene	1.0	U
95-47-6-----	o-Xylene	0.50	U
100-42-5-----	Styrene	0.50	U <i>J</i>

FORM I VOA

FORM 1
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

Lab Name: COMPUCHEM	Method: 8260B	EFLUENT
Lab Code: LIBRTY	Case No.:	SDG No.: 11222
Matrix: (soil/water) WATER		Lab Sample ID: 1122201
Sample wt/vol: 25	(g/ml) ML	Lab File ID: 1122201RA61
Level: (low/med)	LOW	Date Received: 10/12/06
% Moisture: not dec.		Date Analyzed: 10/17/06
GC Column: RTX-VMS	ID: 0.18 (mm)	Dilution Factor: 1.0
Soil Extract Volume: _____	(uL)	Soil Aliquot Volume: _____ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) UG/L	Q
75-25-2-----	Bromoform	0.50	U <i>UJ</i>
79-34-5-----	1,1,2,2-Tetrachloroethane	0.50	U
541-73-1-----	1,3-Dichlorobenzene	0.50	U
106-46-7-----	1,4-Dichlorobenzene	0.50	U
95-50-1-----	1,2-Dichlorobenzene	0.50	U
120-82-1-----	1,2,4-Trichlorobenzene	0.50	U
540-59-0-----	1,2-Dichloroethene (total)	0.63	<i>J</i>
1330-20-7-----	Xylene (total)	0.50	U <i>UJ</i>

FORM I VOA

K1114104

FORM 1
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

Lab Name:	COMPUCHEM	Method:	8270C	EFFLUENT
Lab Code:	LIBRTY	Case No.:	SAS No.:	SDG No.: 11222
Matrix:	(soil/water)	WATER	Lab Sample ID:	1122201
Sample wt/vol:	1000	(g/mL) ML	Lab File ID:	1122201A66
Level:	(low/med)	LOW	Date Received:	10/12/06
% Moisture:	_____	decanted: (Y/N) _____	Date Extracted:	10/17/06
Concentrated Extract Volume:	1000	(uL)	Date Analyzed:	10/19/06
Injection Volume:	1.0	(uL)	Dilution Factor:	1.0
GPC Cleanup:	(Y/N)	N	pH:	_____

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg)	UG/L	Q
111-44-4-----	Bis(2-chloroethyl)ether _____		10	U
106-44-5-----	4-Methylphenol _____		20	U
78-59-1-----	Isophorone _____		10	U
117-81-7-----	bis(2-ethylhexyl)Phthalate _____		10	U

FORM I SV

8270C

11/14/06

FORM 1
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

EFFLUENT

Lab Name: COMPUCHEM Method: 8270C

Lab Code: LIBRTY Case No.: SAS No.: SDG No.: 11222

Matrix: (soil/water) WATER Lab Sample ID: 1122201

Sample wt/vol: 1000 (g/mL) ML Lab File ID: 1122201A60

Level: (low/med) LOW Date Received: 10/12/06

% Moisture: _____ decanted: (Y/N) _____ Date Extracted: 10/17/06

Concentrated Extract Volume: 1000 (uL) Date Analyzed: 10/25/06

Injection Volume: 1.0 (uL) Dilution Factor: 1.0

GPC Cleanup: (Y/N) N pH: _____

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/L Q

CAS NO.	COMPOUND			
87-86-5-----	Pentachlorophenol	0.28	JB	1.00 u

FORM I SV

8270C

11/14/06

1D
GC EXTRACTABLE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

Lab Name:	COMPUCHEM	Contract:	8082	EFFLUENT
Lab Code:	LIBRTY	Case No.:	SAS No.:	SDG No.: 11222
Matrix:	(soil/water) WATER			Lab Sample ID: 1122201
Sample wt/vol:	1000 (g/mL)	ML		Lab File ID: _____
% Moisture:	_____	decanted: (Y/N)	_____	Date Received: 10/12/06
Extraction:	(SepF/Cont/Sonc) SEPF			Date Extracted: 10/13/06
Concentrated Extract Volume:	2500 (uL)			Date Analyzed: 10/15/06
Injection Volume:	1.0 (uL)			Dilution Factor: 1.0
GPC Cleanup:	(Y/N) N	pH:	_____	Sulfur Cleanup: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg)		Q
		UG/L	Q	
12674-11-2-----	Aroclor-1016	0.47	U	
11104-28-2-----	Aroclor-1221	0.63	U	
11141-16-5-----	Aroclor-1232	0.47	U	
53469-21-9-----	Aroclor-1242	0.31	U	
12672-29-6-----	Aroclor-1248	0.31	U	
11097-69-1-----	Aroclor-1254	0.31	U	
11096-82-5-----	Aroclor-1260	0.47	U	

FORM I PEST

SW846 - METALS

-1-

INORGANIC ANALYSES DATA SHEET

EPA SAMPLE NO.

EFFLUENT

Lab Name: COMPUCHEM Contract: _____
 Lab Code: LIBRTY Case No.: _____ SAS No.: _____ SDG No.: 11122
 Matrix (soil/water): WATER Lab Sample ID: 1122201
 Level (low/med): LOW Date Received: 10/12/2006
 % Solids: 0.0

Concentration Units (ug/L or mg/kg dry weight): UG/L

CAS No.	Analyte	Concentration	C	Q	M
7440-38-2	Arsenic	9.5	B		P
7440-41-7	Beryllium	0.45	B	UB	P
7440-43-9	Cadmium	0.20	U		P
7439-97-6	Mercury	0.10	U		CV
7439-96-5	Manganese	0.10	U		P
7782-49-2	Selenium	2.5	U		P
7440-28-0	Thallium	3.2	U		P
7440-66-6	Zinc	1.1	B	UB	P

Color Before: COLORLESS Clarity Before: CLEAR Texture: _____Color After: COLORLESS Clarity After: CLEAR Artifacts: _____Comments: _____



CompuChem a Division of Liberty Analytical Corp.

Remit to: P.O. Box 4603

Cary, NC 27519-4603

Phone: (919) 379-4100

Fax: (919) 379-4050

ANALYTICAL RESULTS

Project: 11222

Project ID: ACS 7010311

Solid results are reported on a dry weight basis.

Lab ID: 1122201 Date Collected: 10/11/2006 02:00 Matrix: Water
Sample ID: EFFLUENT Date Received: 10/12/2006 11:52

Parameters Results Units Report Limit DF Prepared By Analyzed By CAS No. Qual Regl.mt

PH OF WATER 150.1 Analytical Method: EPA 150.1

Analytical Method: EPA 150.1

PH-150.1 7.45 PH 0.00 1 10/19/2006 2477
UNITS

TTL SSPND SOLIDS (TSS) 160.2 W Analytical Method: EPA 160.2

Analytical Method: EPA 160.2

TSS 0.800B mg/L 1.00 1 10/17/2006 2477

J

Date: 10/26/2006

Page 4 of 9

REPORT OF LABORATORY ANALYSIS

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**November 1, 2006 Compliance Sample
Laboratory Results**

FORM 1
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

Lab Name: COMPUCHEM

Method: 8260B

EFFLUENT

Lab Code: LIBRTY Case No.:

SAS No.: SDG No.: 11406

Matrix: (soil/water) WATER

Lab Sample ID: 1140601

Sample wt/vol: 25 (g/ml) ML

Lab File ID: 1140601B62

Level: (low/med) LOW

Date Received: 11/02/06

% Moisture: not dec.

Date Analyzed: 11/02/06

GC Column: RTX-VMS ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: (uL)

Soil Aliquot Volume: (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg)	UG/L	Q
74-87-3-----	Chloromethane	0.50	U	
75-01-4-----	Vinyl Chloride	0.50	U	
74-83-9-----	Bromomethane	0.50	U	
75-00-3-----	Chloroethane	0.30	J	
75-35-4-----	1,1-Dichloroethene	0.50	U	
75-15-0-----	Carbon disulfide	0.50	U	
67-64-1-----	Acetone	2.5	U	UJ
75-09-2-----	Methylene Chloride	0.84		
156-60-5-----	trans-1,2-Dichloroethene	0.50	U	
75-34-3-----	1,1-Dichloroethane	0.15	J	
156-59-2-----	cis-1,2-Dichloroethene	0.42	J	
78-93-3-----	2-butanone	2.5	U	
67-56-3-----	Chloroform	0.50	U	
71-55-6-----	1,1,1-Trichloroethane	0.50	U	UJ
56-23-5-----	Carbon Tetrachloride	0.50	U	UJ
71-43-2-----	Benzene	0.50	U	
107-06-2-----	1,2-Dichloroethane	0.50	U	
79-01-6-----	Trichloroethene	0.50	U	
78-87-5-----	1,2-Dichloropropane	0.50	U	
75-27-4-----	Bromodichloromethane	0.50	U	
10061-01-5-----	cis-1,3-Dichloropropene	0.50	U	
108-10-1-----	4-Methyl-2-pentanone	2.5	U	
108-88-3-----	Toluene	0.50	U	
10061-02-6-----	trans-1,3-Dichloropropene	0.50	U	
79-00-5-----	1,1,2-Trichloroethane	0.50	U	
127-18-4-----	Tetrachloroethene	0.50	U	
591-78-6-----	2-hexanone	2.5	U	
124-48-1-----	Dibromochloromethane	0.50	U	
108-90-7-----	Chlorobenzene	0.50	U	
100-41-4-----	Ethylbenzene	0.50	U	
108-38-3-----	m,p-Xylene	1.0		
95-47-6-----	o-Xylene	0.50	U	
100-42-5-----	Styrene	0.50	U	

FORM I VOA

FORM 1
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

Lab Name: COMPUCHEM

Method: 8260B

EFFLUENT

Lab Code: LIBRTY Case No.:

SAS No.: SDG No.: 11406

Matrix: (soil/water) WATER

Lab Sample ID: 1140601

Sample wt/vol: 25 (g/ml) ML

Lab File ID: 1140601B62

Level: (low/med) LOW

Date Received: 11/02/06

% Moisture: not dec. _____

Date Analyzed: 11/02/06

GC Column: RTX-VMS ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: _____ (uL)

Soil Aliquot Volume: _____ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) UG/L	Q
75-25-2-----	Bromoform	0.50	U
79-34-5-----	1,1,2,2-Tetrachloroethane	0.50	U
541-73-1-----	1,3-Dichlorobenzene	0.50	U
106-46-7-----	1,4-Dichlorobenzene	0.50	U
95-50-1-----	1,2-Dichlorobenzene	0.50	U
120-82-1-----	1,2,4-Trichlorobenzene	0.50	U
540-59-0-----	1,2-Dichloroethene (total)	0.43	J
1330-20-7-----	Xylene (total)	0.50	U

FORM I VOA



COMPUCHEM
a division of Liberty Analytical Corp.

CompuChem a Division of Liberty Analytical Corp.

Remit to: P.O. Box 4603

Cary, NC 27519-4603

Phone: (919) 379-4100

Fax: (919) 379-4050

ANALYTICAL RESULTS

Project: 11406

Project ID: ACS 7010311

Solid results are reported on a dry weight basis.

Lab ID:	1140601	Date Collected:	11/1/2006 00:00	Matrix:	Water
Sample ID:	EFFLUENT	Date Received:	11/2/2006 09:33		
Parameters	Results	Units	Report Limit	DF Prepared	By
				Analyzed	By
				CAS No.	Qual
				RegLmt	
PH OF WATER 150.1			Analytical Method: EPA 150.1		
PH-150.1	7.05	PH UNITS	0.00	1	11/3/2006 2477

Date: 11/08/2006

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REPORT OF LABORATORY ANALYSIS

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**December 21, 2006 Compliance Sample
Laboratory Results**

APPENDIX B

THERMAL OXIDIZER OFF-GAS ANALYTICAL DATA

October 19, 2006 Off-Gas Sample Laboratory Results



AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 1 OFFSITE ISVE

Lab ID#: 0610414A-01A

MODIFIED EPA METHOD TO-14A GC/MS FULL SCAN

File Name:	5110812	Date of Collection:	10/19/06	
Dil. Factor:	1	Date of Analysis:	11/8/06 05:47 PM	
Compound	Rpt. Limit (ppbv)	Amount (ppbv)	Rpt. Limit (uG/m3)	Amount (uG/m3)
Vinyl Chloride	370	330 J	950	840 J
Bromomethane	370	Not Detected	1400	Not Detected
Chloroethane	370	Not Detected	980	Not Detected
1,1-Dichloroethene	370	350 J	1500	1400 J
Methylene Chloride	370	44000	1300	150000
1,1-Dichloroethane	370	4600	1500	19000
cis-1,2-Dichloroethene	370	1400	1500	5600
Chloroform	370	2500	1800	12000
1,1-Trichloroethane	370	36000	2000	190000
Carbon Tetrachloride	370	Not Detected	2300	Not Detected
Benzene	370	18000	1200	57000
1,1-Dichloroethane	370	900	1500	3600
Trichloroethene	370	16000	2000	36000
1,1-Dichloropropane	370	320 J	1700	1500 J
trans-1,3-Dichloropropene	370	Not Detected	1700	Not Detected
Toluene	370	120000	1400	440000
trans-1,3-Dichloropropene	370	Not Detected	1700	Not Detected
1,1,2-Trichloroethane	370	Not Detected	2000	Not Detected
Trichloroethene	370	26000	2500	180000
Chlorobenzene	370	Not Detected	1700	Not Detected
Ethyl Benzene	370	13000	1600	56000
m,p-Xylene	370	50000	1600	220000
p-Xylene	370	18000	1600	80000
Styrene	370	Not Detected	1600	Not Detected
1,1,2-Tetrachloroethane	370	Not Detected	2500	Not Detected
Bromodichloromethane	370	Not Detected	2500	Not Detected
Dibromochloromethane	370	Not Detected	3200	Not Detected
Chloromethane	1500	Not Detected	3100	Not Detected
Acetone	1500	24000	3500	57000
Carbon Disulfide	1500	180 J	4600	550 J
trans-1,2-Dichloroethene	1500	Not Detected	5900	Not Detected
2-Butanone (Methyl Ethyl Ketone)	1500	16000	4400	47000
4-Methyl-2-pentanone	1500	9500	6100	39000
2-Hexanone	1500	280 J	6100	1200 J
Bromoform	1500	Not Detected	15000	Not Detected

- Estimated value.

Container Type: 6 Liter Summa Canister

11/17/06



AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 1 OFFSITE ISVE

Lab ID#: 0610414A-01A

MODIFIED EPA METHOD TO-14A GC/MS FULL SCAN

File Name:	51108123	Date of Collection:	10/19/06
DIL Factor:	741	Date of Analysis:	11/8/06 05:47 PM

Surrogates	%Recovery	Method Limits
1,2-Dichloroethane-d4	98	70-130
Toluene-d8	102	70-130
4-Bromof uorobenzene	114	70-130



AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 2 SBPA ISVE

Lab ID#: 0610414A-02A

MODIFIED EPA METHOD TO-14A GC/MS FULL SCAN

File Name:	5110309	Date of Collection:	10/19/06	
Dil. Factor:	278	Date of Analysis:	11/3/06 06:00 PM	
Compound	Rpt. Limit (ppbv)	Amount (ppbv)	Rpt. Limit (uG/m3)	Amount (uG/m3)
Vinyl Chloride	140	1400	360	3700
Bromomethane	140	Not Detected	540	Not Detected
Chloroethane	140	160	370	420
1,1-Dichloroethene	140	230	550	910
Methylene Chloride	140	4800	480	16000
1,1-Dichloroethane	140	1400	560	5500
cis-1,2-Dichloroethene	140	8700	550	35000
Chloroform	140	3600	680	17000
1,1-Trichloroethane	140	9700	760	53000
Carbon Tetrachloride	140	Not Detected	870	Not Detected
Benzene	140	4700	440	15000
1,1-Dichloroethane	140	240	560	960
Trichloroethene	140	9300	750	50000
1,1-Dichloropropane	140	210	640	970
trans-1,3-Dichloropropene	140	Not Detected	630	Not Detected
Toluene	140	35000	520	130000
trans-1,3-Dichloropropene	140	Not Detected	630	Not Detected
cis-1,2-Trichloroethane	140	Not Detected	760	Not Detected
1,1-Dichloroethene	140	24000	940	160000
Chlorobenzene	140	59 J	640	270 J
Ethyl Benzene	140	8100	600	35000
m,p-Xylene	140	33000	600	140000
p-Xylene	140	12000	600	55000
o-Xylene	140	Not Detected	590	Not Detected
1,1,2,2-Tetrachloroethane	140	Not Detected	950	Not Detected
Bromochloromethane	140	Not Detected	930	Not Detected
Dibromochloromethane	140	Not Detected	1200	Not Detected
Unbromomethane	560	Not Detected	1100	Not Detected
Acetone	560	750	1300	1800
Carbon Disulfide	560	Not Detected	1700	Not Detected
trans-1,2-Dichloroethene	560	Not Detected	2200	Not Detected
2-Butanone (Methyl Ethyl Ketone)	560	420 J	1600	1200 J
4-Methyl-2-pentanone	560	1700	2300	6900
2-Hexanone	560	Not Detected	2300	Not Detected
Bromoform	560	Not Detected	5700	Not Detected

(J = Estimated value.)

Container Type: 6 Liter Summa Canister

OPE
11/17/06



AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 2 SBPA ISVE

Lab ID#: 0610414A-02A

MODIFIED EPA METHOD TO-14A GC/MS FULL SCAN

File Name:	5110309	Date of Collection:	10/19/06
DIL Factor:	278	Date of Analysis:	11/06/06 00:00 PM

Surrogates	%Recovery	Method Limits
1,1-Dichloroethane-d4	98	70-130
Toluene-d8	105	70-130
4-Bromofluorobenzene	116	70-130

ME
11/17/06



AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 3 TOX 1 INF

Lab ID#: 0610414A-03A

MODIFIED EPA METHOD TO-14A GC/MS FULL SCAN

File Name:	5110310	Date of Collection:	10/19/06	
Dil. Factor:	292	Date of Analysis:	11/3/06 08:28 PM	
Compound	Rpt. Limit (ppbv)	Amount (ppbv)	Rpt. Limit (uG/m3)	Amount (uG/m3)
Vinyl Chloride	150	1500	370	3900
Bromomethane	150	Not Detected	570	Not Detected
Chloroethane	150	170	380	450
1,1-Dichloroethene	150	240	580	950
Methylene Chloride	150	4700	510	16000
1,1-Dichloroethane	150	1400	590	5800
cis-1,2-Dichloroethene	150	9000	580	36000
Chloroform	150	3500	710	17000
1,1,1-Trichloroethane	150	10000	800	56000
Carbon Tetrachloride	150	Not Detected	920	Not Detected
Benzene	150	4300	470	14000
1,2-Dichloroethane	150	210	590	870
Trichloroethene	150	9100	780	49000
1,2-Dichloropropane	150	230	670	1000
cis-1,3-Dichloropropene	150	Not Detected	660	Not Detected
toluene	150	34000	550	130000
trans-1,3-Dichloropropene	150	Not Detected	660	Not Detected
1,1,2-Trichloroethane	150	Not Detected	800	Not Detected
Tetrachloroethene	150	26000	990	170000
Chlorobenzene	150	Not Detected	670	Not Detected
Ethyl Benzene	150	8300	630	36000
m-p-Xylene	150	33000	630	140000
o-Xylene	150	13000	630	56000
Syrene	150	Not Detected	620	Not Detected
1,1,2,2-Tetrachloroethane	150	94 J	1000	350 J
Bromodichloromethane	150	Not Detected	980	Not Detected
Dibromochloromethane	150	Not Detected	1200	Not Detected
Chloromethane	580	Not Detected	1200	Not Detected
Acetone	580	800	1400	1900
Carbon Disulfide	580	Not Detected	1800	Not Detected
trans-1,2-Dichloroethene	580	170 J	2300	670 J
2-Butanone (Methyl Ethyl Ketone)	580	410 J	1700	1200 J
4-Methyl-2-pentanone	580	1400	2400	5800
2-Hexanone	580	Not Detected	2400	Not Detected
Chloroform	580	Not Detected	6000	Not Detected

J = Estimated value.

Container Type: 6 Liter Summa Canister

OMS
11/17/06



AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 3 TOX 1 INF

Lab ID#: 0610414A-03A

MODIFIED EPA METHOD TO-14A GC/MS FULL SCAN

File Name:	3TOX1INF.DAT	Date of Collection:	10/19/06
DIL. Factor:	292	Date of Analysis:	11/3/06 06:28 PM

Surrogates	%Recovery	Method Limits
1,2-Dichloroethane-d4	98	70-130
Toluene-d8	100	70-130
4-Hromof uorobenzene	116	70-130

11/17/06
JAS



AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 4 TOX 1 INF DUP

Lab ID#: 0610414A-04A

MODIFIED EPA METHOD TO-14A GC/MS FULL SCAN

File Name	5110311	Date of Collection	10/19/06	
Dil. Factor	272	Date of Analysis	11/3/06 06:55 PM	
Compound	Rpt. Limit (ppbv)	Amount (ppbv)	Rpt. Limit (uG/m3)	Amount (uG/m3)
Vinyl Chloride	140	1800 R	350	4500
Bromomethane	140	Not Detected	530	Not Detected
Chloroethane	140	220	360	570
1,1-Dichloroethene	140	220	540	880
Methylene Chloride	140	5000	470	17000
1,1-Dichloroethane	140	1400	550	5700
cis-1,2-Dichloroethene	140	9700	540	38000
Chloroform	140	3400	660	17000
1,1,1-Trichloroethane	140	10000	740	53000
Carbon Tetrachloride	140	Not Detected	860	Not Detected
Benzene	140	4000	430	13000
1,2-Dichloroethane	140	220	550	880
Trichloroethene	140	9300	730	50000
1,2-Dichloropropane	140	190	630	880
cis-1,3-Dichloropropene	140	Not Detected	620	Not Detected
Toluene	140	34000	510	130000
trans-1,3-Dichloropropene	140	Not Detected	620	Not Detected
1,2-Trichloroethane	140	Not Detected	740	Not Detected
Tetrachloroethene	140	26000	920	180000
Chlorobenzene	140	58 J	630	260 J
Methyl Benzene	140	8000	590	35000
m,p-Xylene	140	32000	590	140000
o-Xylene	140	13000	590	55000
Styrene	140	Not Detected	580	Not Detected
1,1,2,2-Tetrachloroethane	140	Not Detected	930	Not Detected
Bromodichromethane	140	Not Detected	910	Not Detected
Dibromochromomethane	140	Not Detected	1200	Not Detected
Chloromethane	540	Not Detected	1100	Not Detected
Acetone	540	620	1300	1500
Carbon Disulfide	540	Not Detected	1700	Not Detected
trans-1,2-Dichloroethene	540	110 J	2200	430 J
2-Butanone (Methyl Ethyl Ketone)	540	390 J	1600	1100 J
1-Methyl-2-pentanone	540	1600	2200	6500
2-Hexanone	540	Not Detected	2200	Not Detected
Bromoform	540	Not Detected	5600	Not Detected

J = Estimated value.

Container Type: 6 Liter Summa Canister

11/17/06



AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 4 TOX 1 INF DUP

Lab ID#: 0610414A-04A

MODIFIED EPA METHOD TO-14A GC/MS FULL SCAN

File Name:	4TOX1INF.DAT	Date of Collection:	10/19/06
DIL Factor:	272	Data of Analysis:	11/3/06 06:55 PM

Surrogates	%Recovery	Method Limits
1,2-Dichloroethane-d4	97	70-130
Toluene-d8	101	70-130
4-Isomeric uorobenzene	117	70-130



AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 5 TOX 1 EFF

Lab ID#: 0610414A-05A

MODIFIED EPA METHOD TO-14A GC/MS FULL SCAN

File Name	5110312	Date of Collection	10/13/06	
Dil Factor	1.39	Date of Analysis	11/3/06 07:27 PM	
Compound	Rpt. Limit (ppbv)	Amount (ppbv)	Rpt. Limit (uG/m3)	Amount (uG/m3)
Vinyl Chloride	0.70	30 R	1.8	76
Bromomethane	0.70	Not Detected	2.7	Not Detected
Chloroethane	0.70	1.9	1.8	5.0
1,1-Dichloroethene	0.70	57	2.8	220
Methylene Chloride	0.70	170	2.4	600
1,1-Dichloroethane	0.70	5.3	2.8	22
cis-1,2-Dichloroethene	0.70	58	2.8	230
Chloroform	0.70	16	3.4	80
1,1,1-Trichloroethane	0.70	35	3.8	190
Carbon Tetrachloride	0.70	0.55 J	4.4	3.4 J
Benzene	0.70	72	2.2	230
,,Dichloroethane	0.70	1.1	2.8	4.5
Trichloroethene	0.70	72	3.7	380
1,1-Dichloropropane	0.70	0.76	3.2	3.5
trans-1,3-Dichloropropene	0.70	0.21 J	3.2	0.96 J
Toluene	0.70	130	2.6	490
trans-1,3-Dichloropropene	0.70	Not Detected	3.2	Not Detected
,,2-Trichloroethane	0.70	Not Detected	3.8	Not Detected
1,1,1-Tetrachloroethene	0.70	220	4.7	1400
Chlorobenzene	0.70	1.9	3.2	8.8
Ethyl Benzene	0.70	32	3.0	140
m,p-Xylene	0.70	130	3.0	560
,,Xylene	0.70	54	3.0	230
Syrene	0.70	17	3.0	72
1,1,2,2-Tetrachloroethane	0.70	0.76	4.8	5.2
Bromodichloromethane	0.70	Not Detected	4.6	Not Detected
D bromochloromethane	0.70	Not Detected	5.9	Not Detected
Chloromethane	2.8	9.0	5.7	18
Acetone	2.8	85	6.6	200
Carbon Disulfide	2.8	1.2 J	8.6	3.7 J
trans-1,2-Dichloroethene	2.8	14	11	55
2-Butanone (Methyl Ethyl Ketone)	2.8	14	8.2	42
4-Methyl -2-pentanone	2.8	17	11	70
,, Hexanone	2.8	0.69 J	11	2.8 J
Bromoform	2.8	Not Detected	29	Not Detected

J = Estimated value.

Container Type: 6 Liter Summa Canister

CNG
11/17/06



AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 5 TOX 1 EFF

Lab ID#: 0610414A-05A

MODIFIED EPA METHOD TO-14A GC/MS FULL SCAN

File Name:	5110312	Date of Collection:	10/19/08
DIL. Factor:	1.39	Date of Analysis:	11/3/08 07:27 PM

Surrogates	%Recovery	Method Limits
1,2-Dichloroethane-d4	104	70-130
Toluene-d8	102	70-130
4-Bromofluorobenzene	114	70-130



AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 6 TOX 2 INF

Lab ID#: 0610414A-06A

MODIFIED EPA METHOD TO-14A GC/MS FULL SCAN

File Name:	5110813	Date of Collection:	10/19/06	
Dil Factor:	74	Date of Analysis:	11/8/06 06:15 PM	
Compound	Rpt. Limit (ppbv)	Amount (ppbv)	Rpt. Limit (uG/m3)	Amount (uG/m3)
Vinyl Chloride	370	1500	950	3800
Bromomethane	370	Not Detected	1400	Not Detected
Chloroethane	370	570	980	1500
1,1-Dichloroethene	370	300 J	1500	1200 J
Methylene Chloride	370	30000	1300	110000
1,1-Dichloroethane	370	3800	1500	15000
1,1,2,2-Tetrachloroethene	370	6000	1500	24000
Chloroform	370	1800	1800	3900
1,1,1-Trichloroethane	370	29000	2000	160000
Carbon Tetrachloride	370	Not Detected	2300	Not Detected
Benzene	370	15000	1200	48000
1,2-Dichloroethane	370	820	1500	3300
Trichloroethene	370	13000	2000	70000
1,2-Dichloropropane	370	290 J	1700	1300 J
cis-1,3-Dichloropropene	370	Not Detected	1700	Not Detected
Toluene	370	89000	1400	340000
trans-1,3-Dichloropropene	370	Not Detected	1700	Not Detected
1,1,2-Trichloroethane	370	Not Detected	2000	Not Detected
Tetrachloroethene	370	23000	2500	160000
Chlorobenzene	370	Not Detected	1700	Not Detected
Ethyl Benzene	370	9800	1600	42000
m,p-Xylene	370	37000	1600	160000
c-Xylene	370	13000	1600	56000
Stryrene	370	Not Detected	1600	Not Detected
1,1,2,2-Tetrachloroethane	370	Not Detected	2500	Not Detected
Bromodichloromethane	370	Not Detected	2500	Not Detected
D bromochloromethane	370	Not Detected	3200	Not Detected
Chloromethane	1500	230 J	3100	480 J
Acetone	1500	16000	3500	37000
Carbon Disulfide	1500	210 J	4600	650 J
trans-1,2-Dichloroethene	1500	Not Detected	5900	Not Detected
2 Butanone (Methyl Ethyl Ketone)	1500	12000	4400	35000
4 Methyl-2-pentanone	1500	6600	6100	27000
2 Hexanone	1500	Not Detected	6100	Not Detected
Bromoform	1500	Not Detected	15000	Not Detected

J = Estimated value.

Container Type: 6 Liter Summa Canister

CJS
11/17/06



AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 6 TOX 2 INF

Lab ID#: 0610414A-06A

MODIFIED EPA METHOD TO-14A GC/MS FULL SCAN

File Name:	5110813	Date of Collection:	10/19/06
DIL Factor:	1/741	Data of Analysis:	11/06/06 06:15 PM

Surrogates	%Recovery	Method Limits
1,2-Dichloroethane-d4	98	70-130
Toluene-d8	99	70-130
4-Bromofluorobenzene	117	70-130

CRS
11/17/06



AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 7 TOX 2 INF DUP

Lab ID#: 0610414A-07A

MODIFIED EPA METHOD TO-14A GC/MS FULL SCAN

File Name	5110314	Date of Collection	10/19/06	
Dil. Factor	644	Date of Analysis	11/3/06 08:23 PM	
Compound	Rpt. Limit (ppbv)	Amount (ppbv)	Rpt. Limit (uG/m3)	Amount (uG/m3)
Vinyl Chloride	320	1400	820	3500
Bromomethane	320	Not Detected	1200	Not Detected
Chloroethane	320	540	850	1400
1,1-Dichloroethene	320	350	1300	1400
Methylene Chloride	320	30000	1100	100000
1,1-Dichloroethane	320	3600	1300	14000
cis-1,2-Dichloroethene	320	6800	1300	27000
Chloroform	320	1900	1600	9100
1,1,1-Trichloroethane	320	26000	1800	140000
Carbon Tetrachloride	320	Not Detected	2000	Not Detected
Benzene	320	15000	1000	48000
1,2-Dichloroethane	320	780	1300	3200
Trichloroethene	320	14000	1700	78000
1,2-Dichloropropane	320	240 J	1500	1100 J
cis-1,3-D chloropropene	320	Not Detected	1500	Not Detected
Toluene	320	100000	1200	380000
trans-1,3-Dichloropropene	320	Not Detected	1500	Not Detected
1,1,2-Trichloroethane	320	Not Detected	1800	Not Detected
Tetrachloroethene	320	24000	2200	160000
Chlorobenzene	320	Not Detected	1500	Not Detected
Ethyl Benzene	320	12000	1400	53000
m,p-Xylene	320	51000	1400	220000
c-Xylene	320	18000	1400	76000
Syrene	320	Not Detected	1400	Not Detected
1,1,2,2-Tetrachloroethane	320	Not Detected	2200	Not Detected
Bromodichloromethane	320	Not Detected	2200	Not Detected
Dibromochloromethane	320	Not Detected	2700	Not Detected
Chloromethane	1300	Not Detected	2600	Not Detected
Acetone	1300	19000	3000	44000
Carbon Disulfide	1300	Not Detected	4000	Not Detected
trans-1,2-Dichloroethene	1300	Not Detected	5100	Not Detected
2-Butanone (Methyl Ethyl Ketone)	1300	12000	3800	36000
4-Methyl-2-pentanone	1300	6100	5300	25000
2-Hexanone	1300	Not Detected	5300	Not Detected
Bromocform	1300	Not Detected	13000	Not Detected

J = Estimated value.

Container Type: 6 Liter Summa Canister

11/17/06



AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 7 TOX 2 INF DUP

Lab ID#: 0610414A-07A

MODIFIED EPA METHOD TO-14A GC/MS FULL SCAN

File Name:	7_TOX_2_INF_DUP_5110314	Date of Collection:	10/19/06
Dil. Factor:	1/1	Date of Analysis:	11/3/06 08:23 PM

Surrogates	%Recovery	Method Limits
1,2-Dichloroethane-d4	96	70-130
Toluene-d8	102	70-130
4-Bromof uorobenzene	113	70-130

CHL
11/17/06



AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 8 TOX 2 EFF

Lab ID#: 0610414A-08A

MODIFIED EPA METHOD TO-14A GC/MS FULL SCAN

File Name:	8TOX2EFF	Date of Collection:	10/19/06	
Dil. Factor:	11.1	Date of Analysis:	11/3/06 08:50 PM	
Compound	Rpt. Limit (ppbv)	Amount (ppbv)	Rpt. Limit (uG/m3)	Amount (uG/m3)
Vinyl Chloride	5.6	57	14	150
Bromomethane	5.6	Not Detected	22	Not Detected
Chloroethane	5.6	10	15	27
1,1-Dichloroethene	5.6	160	22	620
Methylene Chloride	5.6	580	19	2000
1,1-Dichloroethane	5.6	60	22	240
cis-1,2-Dichloroethene	5.6	120	22	480
Chloroform	5.6	36	27	180
1,1,1-Trichloroethane	5.6	430	30	2300
Carbon Tetrachloride	5.6	3.8 J	35	24 J
Benzene	5.6	480	18	1500
1,2-Dichloroethane	5.6	13	22	52
Trichloroethene	5.6	290	30	1600
1,2-Dichloropropane	5.6	3.8 J	26	18 J
trans-1,3-Dichloropropene	5.6	Not Detected	25	Not Detected
Toluene	5.6	1400	21	5400
trans-1,3-Dichloropropene	5.6	Not Detected	25	Not Detected
1,1,2-Trichloroethane	5.6	Not Detected	30	Not Detected
Tetrachloroethene	5.6	570	38	3900
Chlorobenzene	5.6	4.0 J	26	18 J
Ethyl Benzene	5.6	120	24	540
m,p-Xylene	5.6	430	24	1900
c-Xylene	5.6	150	24	670
Styrene	5.6	95	24	410
1,1,2,2-Tetrachloroethane	5.6	Not Detected	38	Not Detected
Bromodichloromethane	5.6	Not Detected	37	Not Detected
D bromochloromethane	5.6	Not Detected	47	Not Detected
Chloromethane	22	Not Detected	46	Not Detected
Acetone	22	470	53	1100
Carbon Disulfide	22	9.2 J	69	29 J
trans-1,2-Dichloroethene	22	22	88	88
2 Butanone (Methyl Ethyl Ketone)	22	200	65	590
4-Methyl-2-pentanone	22	61	91	250
2-Hexanone	22	2.6 J	91	10 J
Bromoform	22	Not Detected	230	Not Detected

J = Estimated value.

Container Type: 6 Liter Summa Canister

CHL
11/17/06



AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 8 TOX 2 EFF

Lab ID#: 0610414A-08A

MODIFIED EPA METHOD TO-14A GC/MS FULL SCAN

File Name:	5110315	Date of Collection:	10/19/06
DIL Factor:	11.1	Date of Analysis:	11/3/06 08:50 PM

Surrogates	%Recovery	Method Limits
1,1-Dichloroethane-d4	94	70-130
Toluene-d8	103	70-130
4-Bromofluorobenzene	113	70-130



AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 1 OFFSITE ISVE

Lab ID#: 0610414B-01A

MODIFIED EPA METHOD TO-13A GC/MS FULL SCAN

File Name:	ISVE_0610414B-01A	Date of Collection:	10/19/06
Dil. Factor:	1.00	Date of Analysis:	10/23/06 12:47 PM
			Date of Extraction:
			10/20/06

Compound	Rpt. Limit (ug)	Amount (ug)
Phenol	5.0	Not Detected
bis(2-Chloroethyl) Ether	1.0	9.2
2-Chlorophenol	5.0	Not Detected
1,3-Dichlorobenzene	1.0	2.6
1,4-Dichlorobenzene	1.0	10
1,2-Dichlorobenzene	1.0	88
2-Methylphenol (o-Cresol)	5.0	Not Detected
N-Nitroso-di-n-propylamine	1.0	Not Detected
4-Methylphenol/3-Methylphenol	5.0	4.8 J 15
Hexachlorcethane	1.0	Not Detected
Nitrobenzene	1.0	Not Detected
Isophorone	1.0	49
2-Nitrophenol	5.0	Not Detected
2,4-D-methylphenol	5.0	Not Detected
bis(2-Chloroethoxy) Methane	1.0	Not Detected
2,4-Dichlorophenol	5.0	Not Detected
1,2,4-Trichlorobenzene	1.0	4.5
Naphthalene	1.0	100
4-Chloraniline	10	Not Detected
Hexachlorobutadiene	1.0	5.9
4-Chloro-3-methylphenol	5.0	Not Detected
2-Methylnaphthalene	1.0	21
Hexachlorocyclopentadiene	20	Not Detected
2,4,6-Trichlorophenol	5.0	Not Detected
2,4,5-Trichlorophenol	5.0	Not Detected
2-Chloronaphthalene	1.0	Not Detected
2-Nitroaniline	10	Not Detected
Dimethylphthalate	5.0	Not Detected
Acenaphthylene	1.0	Not Detected
2,6-Dinitrotoluene	5.0	Not Detected
3-Nitroaniline	10	Not Detected
Acenaphthene	1.0	Not Detected
2,4-Dinitrophenol	20	Not Detected
4-Nitrophenol	20	Not Detected
2,4-Dinitrotoluene	5.0	Not Detected
Benzofuran	1.0	Not Detected
Diethylphthalate	5.0	Not Detected
Fluorene	1.0	Not Detected

09
11/7/06



AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 1 OFFSITE ISVE

Lab ID#: 0610414B-01A

MODIFIED EPA METHOD TO-13A GC/MS FULL SCAN

File Name:	0610414B-01A	Date of Collection:	10/19/06	
DIL Factor:	1.00	Date of Analysis:	10/23/06 12:47 PM	
			Date of Extraction:	10/20/06

Compound	Rpt. Limit (μ g)	Amount (μ g)
4-Chlorophenyl-phenyl Ether	1.0	Not Detected
4-Nitroaniline	10	Not Detected
4,6-Dinitro-2-methylphenol	10	Not Detected
N-Nitrosodiphenylamine	10	Not Detected
4-Bromophenyl-phenyl Ether	1.0	Not Detected
Heptachlorobenzene	1.0	Not Detected
Pentachlorophenol	20	Not Detected
Phenanthrene	1.0	Not Detected
Anthracene	1.0	Not Detected
di-n-Butylphthalate	5.0	0.99 J 15
Fluoranthene	1.0	Not Detected
Pyrene	1.0	Not Detected
Ethylbenzylphthalate	5.0	Not Detected
3,3'-Dichlorobenzidine	20	Not Detected
Chrysene	1.0	Not Detected
Benzo(a)anthracene	1.0	Not Detected
Is(2-Ethylhexyl)phthalate	5.0	4.2 J 15
Di-n-Octylphthalate	5.0	Not Detected
Benzo(b)fluoranthene	1.0	Not Detected
Benzo(k)fluoranthene	1.0	Not Detected
Benzo(a)pyrene	1.0	Not Detected
Indeno(1,2,3-c,d)pyrene	1.0	Not Detected
Dibenz(a,h)anthracene	1.0	Not Detected
Benzo(g,h,i)perylene	1.0	Not Detected

J = Estimated value.

Container Type: XAD Tube

Surrogates	%Recovery	Method Limits
2-Fluoropheno	61	50-150
Phenol-c5	83	50-150
Nitrobenzene-c5	95	50-150
2,4,6-Tribromophenol	71	50-150
Fluorene-d10	75	60-120
Pyrene-c10	77	60-120



AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 2 SBPA ISVE

Lab ID#: 0610414B-02A

MODIFIED EPA METHOD TO-13A GC/MS FULL SCAN

File Name:	D102312	Date of Collection:	10/19/06
DIL Factor:	100	Date of Analysis:	10/23/06 01:17 PM
			Time of Extraction: 07/2006

Compound	Rpt. Limit (ug)	Amount (ug)
Phenol	5.0	Not Detected
bis(2-Chloroethyl) Ether	1.0	2.1
2-Chlorophenol	5.0	Not Detected
1,3-Dichlorobenzene	1.0	2.3
1,4-Dichlorobenzene	1.0	5.0
1,2-Dichlorobenzene	1.0	22
2-Methylphenol (o-Cresol)	5.0	Not Detected
N-Nitroso-di-n-propylamine	1.0	Not Detected
4-Methylphenol/3-Methylphenol	5.0	Not Detected
Hexachloroethane	1.0	Not Detected
Nitrobenzene	1.0	Not Detected
Isophorone	1.0	5.9
2-Nitrophenol	5.0	Not Detected
2,4-Dimethylphenol	5.0	Not Detected
bis(2-Chloroethoxy) Methane	1.0	Not Detected
2,4-Dichlorophenol	5.0	Not Detected
1,2,4-Trichlorobenzene	1.0	0.53 J 15
Naphthalene	1.0	22
4-Chloroaniline	10	Not Detected
Hexachlorobutadiene	1.0	8.2
4-Chloro-3-methylphenol	5.0	Not Detected
2-Methylnaphthalene	1.0	12
Hexachlorocyclopentadiene	20	Not Detected
2,4,6-Trichlorophenol	5.0	Not Detected
2,4,5-Trichlorophenol	5.0	Not Detected
2-Chloronaphthalene	1.0	Not Detected
2-Nitroaniline	10	Not Detected
Dimethylphthalate	5.0	Not Detected
Acenaphthylene	1.0	Not Detected
2,6-Dinitrotoluene	5.0	Not Detected
3-Nitroaniline	10	Not Detected
Acenaphthene	1.0	Not Detected
2,4-Dinitrophenol	20	Not Detected
4-Nitrophenol	20	Not Detected
2,4-Dinitrotoluene	5.0	Not Detected
D benzofuran	1.0	Not Detected
Dethylphthalate	5.0	1.3 J 15
Fluorene	1.0	Not Detected

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10/10/06



AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 2 SBPA ISVE

Lab ID#: 0610414B-02A

MODIFIED EPA METHOD TO-13A GC/MS FULL SCAN

File Name:	2SBPAISVE.DAT	Date of Collection:	10/19/06
Dil. Factor:	1.00	Date of Analysis:	10/23/06 01:11:21
			Date of Extraction: 10/20/06

Compound	Rpt. Limit (μ g)	Amount (μ g)
4-Chlorophenyl-phenyl Ether	1.0	Not Detected
4-N-trcamiline	10	Not Detected
4,6-Dinitro-2-methylphenol	10	Not Detected
N-Nitrosodiphenylamine	10	Not Detected
4-Bromophenyl-phenyl Ether	1.0	Not Detected
Hexachlorobenzene	1.0	Not Detected
Partachorophenol	20	Not Detected
Phenanthrene	1.0	Not Detected
Anthracene	1.0	Not Detected
d-n-Butylphthalate	5.0	Not Detected
Fluoranthene	1.0	Not Detected
Pyrene	1.0	Not Detected
Butylbenzylphthalate	5.0	Not Detected
3,3'-D chlorobenzidine	20	Not Detected
Chrysene	1.0	Not Detected
Benz(a)anthracene	1.0	Not Detected
b-s(2-Ethylhexyl)phthalate	5.0	14
D-n-Octylphthalate	5.0	Not Detected
Benz(b)fluoranthene	1.0	Not Detected
Benz(k)fluoranthene	1.0	Not Detected
Benz(a)pyrene	1.0	Not Detected
Indeno(1,2,3-c,d)pyrene	1.0	Not Detected
Dibenz(a,h)anthracene	1.0	Not Detected
Benzo(g,h,i)perylene	1.0	Not Detected

J = Estimated value.

Container Type: XAD Tube

Surrogates	%Recovery	Method Limits
2-Fluorophenol	71	50-150
Phenol-d5	74	50-150
Mitrobenzene-d5	76	50-150
2,4,6-Tribromophenol	62	50-150
Fluorene-d ¹⁰	76	60-120
Pyrene-d ¹⁰	77	60-120

OS
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AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 2 SBPA ISVE Duplicate

Lab ID#: 0610414B-02AA

MODIFIED EPA METHOD TO-13A GC/MS FULL SCAN

File Name:	P102313	Date of Collection:	10/19/06
Dil. Factor:	1.00	Date of Analysis:	10/23/06 01:47 PM
		Date of Extraction:	10/23/06

Compound	Rpt. Limit (ug)	Amount (ug)
Phenol	5.0	Not Detected
bis(2-Chloroethyl) Ether	1.0	2.2
2-Chlorophenol	5.0	Not Detected
1,3-Dichlorobenzene	1.0	2.1
1,4-Dichlorobenzene	1.0	5.1
1,2-Dichlorobenzene	1.0	22
2-Methylphenol (o-Cresol)	5.0	Not Detected
N-Nitroso-di-n-propylamine	1.0	Not Detected
4-Methylpheno/3-Methylphenol	5.0	Not Detected
Hexachloroethane	1.0	Not Detected
Nitrobenzene	1.0	Not Detected
Isophorone	1.0	5.9
2-Nitrophenol	5.0	Not Detected
2,4-Dimethylphenol	5.0	Not Detected
bis(2-Chloroethoxy) Methane	1.0	Not Detected
2,4-Dichlorophenol	5.0	Not Detected
1,2,4-Trichlorobenzene	1.0	0.56 J 15
Naphthalene	1.0	22
4-Chloraniline	10	Not Detected
Hexachlorobutadiene	1.0	7.7
4-Chloro-3-methylphenol	5.0	Not Detected
2-Methylnaphthalene	1.0	12
Hexachlorocyclopentadiene	20	Not Detected
2,4,6-Trichlorophenol	5.0	Not Detected
2,4,5-Trichlorophenol	5.0	Not Detected
2-Chloronaphthalene	1.0	Not Detected
2-Nitroaniline	10	Not Detected
C-methylphthalate	5.0	Not Detected
Acenaphthylene	1.0	Not Detected
2,6-Dinitrotoluene	5.0	Not Detected
3-Nitroaniline	10	Not Detected
Acenaphthene	1.0	Not Detected
2,4-Dinitrophenol	20	Not Detected
4-Nitrophenol	20	Not Detected
2,4-Dinitrotoluene	5.0	Not Detected
Dibenzofuran	1.0	Not Detected
Diethylphthalate	5.0	1.3 J 15
Fluorene	1.0	Not Detected

06/11/06



AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 2 SBPA ISVE Duplicate

Lab ID#: 0610414B-02AA

MODIFIED EPA METHOD TO-13A GC/MS FULL SCAN

File Name:	P102313	Date of Collection:	10/18/06	
Dil Factor:	1.00	Date of Analysis:	10/23/06 01:47 PM	
			Date of Extraction:	10/20/06

Compound	Rpt. Limit (μ g)	Amount (μ g)
4-Chlorophenyl-phenyl Ether	1.0	Not Detected
4-Nitroaniline	10	Not Detected
4,6-Dinitro-2-methylphenol	10	Not Detected
N-Nitrosodiphenylamine	10	Not Detected
4-Bromophenyl-phenyl Ether	1.0	Not Detected
Hexachlorobenzene	1.0	Not Detected
Pentachlorophenol	20	Not Detected
Phenanthrene	1.0	Not Detected
Anthracene	1.0	Not Detected
Di-n-Butylphthalate	5.0	Not Detected
Fluoranthene	1.0	Not Detected
Pyrene	1.0	Not Detected
Benzylbenzylphthalate	5.0	Not Detected
3,3'-Dichlorobenzidine	20	Not Detected
Chrysene	1.0	Not Detected
Benz(a)anthracene	1.0	Not Detected
bis(2-Ethylhexyl)phthalate	5.0	14
D- <i>n</i> -Octylphthalate	5.0	Not Detected
Benz(b)fluoranthene	1.0	Not Detected
Benz(k)fluoranthene	1.0	Not Detected
Benz(a)pyrene	1.0	Not Detected
Endo(1,2,3-c,d)pyrene	1.0	Not Detected
Obenz(a,h)anthracene	1.0	Not Detected
Benz(g,h,i)perylene	1.0	Not Detected

J = Estimated value.

Container Type: XAD Tube

Surrogates	%Recovery	Method Limits
2-Fluorophenol	69	50-150
Phenol-d5	75	50-150
Nitrobenzene-d5	75	50-150
2,4,6-Tribromophenol	60	50-150
Fluorene-d10	75	60-120
Pyrene-d10	77	60-120

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AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 3 TOX 1 INF

Lab ID#: 0610414B-03A

MODIFIED EPA METHOD TO-13A GC/MS FULL SCAN

File Name:	p102314	Date of Collection:	10/19/06
Dil Factor:	100	Date of Analysis:	10/23/06 02:17 PM
Sample ID:		Date of Extraction:	10/20/06

Compound	Rpt. Limit (μ g)	Amount (μ g)
Phenol	5.0	Not Detected
bis(2-Chloroethyl) Ether	1.0	2.2
2-Chlorophenol	5.0	Not Detected
1,3-Dichlorobenzene	1.0	2.6
1,4-Dichlorobenzene	1.0	5.7
1,2-Dichlorobenzene	1.0	24
2-Methylphenol (o-Cresol)	5.0	Not Detected
N-Nitroso-di-n-propylamine	1.0	Not Detected
4-Methylphenol/3-Methylphenol	5.0	Not Detected
Hexachloroethane	1.0	Not Detected
Nitrobenzene	1.0	Not Detected
Isophorone	1.0	3.0
2-Nitrophenol	5.0	Not Detected
2,4-Dimethylphenol	5.0	Not Detected
bis(2-Chloroethoxy) Methane	1.0	Not Detected
2,4-Dichlorophenol	5.0	Not Detected
1,2,4-Trichlorobenzene	1.0	0.72 J
Naphthalene	1.0	25
4-Chloroaniline	10	Not Detected
Hexachlorobutadiene	1.0	8.8
4-Chloro-3-methylphenol	5.0	Not Detected
2-Methylnaphthalene	1.0	14
Hexachlorocyclopentadiene	20	Not Detected
2,4,6-Trichlorophenol	5.0	Not Detected
2,4,5-Trichlorophenol	5.0	Not Detected
2-Chloronaphthalene	1.0	Not Detected
2-Nitroaniline	10	Not Detected
Dimethyl phthalate	5.0	Not Detected
Acenaphthylene	1.0	Not Detected
2,6-Dinitrotoluene	5.0	Not Detected
3-Nitroaniline	10	Not Detected
Acenaphthene	1.0	Not Detected
2,4-Dinitrophenol	20	Not Detected
4-Nitrophenol	20	Not Detected
2,4-Dinitrotoluene	5.0	Not Detected
E benzofuran	1.0	Not Detected
E ethylenthalate	5.0	Not Detected
Fluorene	1.0	Not Detected



AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 3 TOX 1 INF

Lab ID#: 0610414B-03A

MODIFIED EPA METHOD TO-13A GC/MS FULL SCAN

File Name	0610414B-03A	Date of Collection	10/19/06
Dil. Factor	1.00	Date of Analysis	10/23/06 02:17 PM
		Date of Extraction	10/20/06

Compound	Rpt. Limit (μ g)	Amount (μ g)
1-Chlorophenyl-phenyl Ether	1.0	Not Detected
1-Nitroaniline	10	Not Detected
4,6-Dinitro-2-methylphenol	10	Not Detected
4-Nitrosodiphenylamine	10	Not Detected
4-Bromophenyl-phenyl Ether	1.0	Not Detected
Hexachlorobenzene	1.0	Not Detected
Heptachlorophenol	20	Not Detected
1-Benanthrene	1.0	Not Detected
1-Phenanthrene	1.0	Not Detected
1-n-Butylphthalate	5.0	Not Detected
Fluoranthene	1.0	Not Detected
1,3-Phe	1.0	Not Detected
1-Benzylphthalate	5.0	Not Detected
1,3'-Dichlorobenzidine	20	Not Detected
Chrysene	1.0	Not Detected
Benzo(a)anthracene	1.0	Not Detected
bis(2-Ethylhexyl)phthalate	5.0	Not Detected
1-n-Octylphthalate	5.0	Not Detected
Benzo(b)fluoranthene	1.0	Not Detected
Benzo(k)fluoranthene	1.0	Not Detected
Benzo(a)pyrene	1.0	Not Detected
1,4-deno(1,2,3-c,d)pyrene	1.0	Not Detected
Dibenz(a,h)anthracene	1.0	Not Detected
Benzo(g,h,i)perylene	1.0	Not Detected

J = Estimated value.

Container Type: XAD Tube

Surrogates	%Recovery	Method Limits
2-Fluorophenol	69	50-150
Pheno-d5	77	50-150
Nitrobenzene-d5	77	50-150
2,4,6-Tribromophenol	60	50-150
Fluorene-d10	76	60-120
Fyrene-d10	77	60-120

CHS
11/17/06



AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 4 TOX 1 INF DUP

Lab ID#: 0610414B-04A

MODIFIED EPA METHOD TO-13A GC/MS FULL SCAN

File Name:	4 TOX 1 INF DUP102320	Date of Collection:	10/19/06
Dil. Factor:	1.00	Date of Analysis:	10/23/06 05:16 PM
		Date of Extraction:	10/20/06

Compound	Rpt. Limit (μ g)	Amount (μ g)
Phenol	5.0	Not Detected
4-(2-Chloroethyl) Ether	1.0	2.8
4-Chlorophenol	5.0	Not Detected
1,3-Dichlorobenzene	1.0	2.9
1,4-Dichlorobenzene	1.0	6.7
1,2-Dichlorobenzene	1.0	28
2-Methylphenol (o-Cresol)	5.0	Not Detected
4-Nitroso-di-n-propylamine	1.0	Not Detected
4-Methylphenol/3-Methylphenol	5.0	Not Detected
Hexachloroethane	1.0	Not Detected
Toluene	1.0	Not Detected
Phenone	1.0	3.8
4-Nitrophenol	5.0	Not Detected
2,4-Dimethylphenol	5.0	Not Detected
2-(2-Chloroethoxy) Methane	1.0	Not Detected
2,3-Dichloropheno	5.0	Not Detected
1,2,4-Trichlorobenzene	1.0	0.91 J
Naphthalene	1.0	32
4-Chloraniline	10	Not Detected
1-Pachlorobutadiene	1.0	11
4-Chloro-3-methylphenol	5.0	Not Detected
2 Methylnaphthalene	1.0	19
Hexachlorocyclopentadiene	20	Not Detected
2,4,6-Trichlorophenol	5.0	Not Detected
2,4,5-Trichlorophenol	5.0	Not Detected
2-Chloronaphtha-ene	1.0	Not Detected
2-Nitroaniline	10	Not Detected
Dimethylphthalate	5.0	Not Detected
Acenaphthylene	1.0	Not Detected
2,6-Dinitrotoluene	5.0	Not Detected
3-Nitroaniline	10	Not Detected
Acenaphthene	1.0	Not Detected
2,4-Dinitropheno	20	Not Detected
4-Nitrophenol	20	Not Detected
2,4-Dinitrotoluene	5.0	Not Detected
Dibenzofuran	1.0	Not Detected
Diethylphthalate	5.0	1.0 J
Clorene	1.0	Not Detected



AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 4 TOX 1 INF DUP

Lab ID#: 0610414B-04A

MODIFIED EPA METHOD TO-13A GC/MS FULL SCAN

File Name:	P102320	Date of Collection:	10/19/06
Dil. Factor:	1.00	Date of Analysis:	10/23/06 05:16 PM
		Date of Extraction:	10/20/06 05:16 PM

Compound	Rpt. Limit (ug)	Amount (ug)
4-Chlorophenyl-phenyl Ether	1.0	Not Detected
4-Nitroaniline	10	Not Detected
4,6-Dinitro-2-methylphenol	10	Not Detected
N-Nitrosodiphenylamine	10	Not Detected
4-Bromophenyl-phenyl Ether	1.0	Not Detected
Hexachlorobenzene	1.0	Not Detected
Pentachlorophenol	20	Not Detected
Phenanthrene	1.0	Not Detected
Anthracene	1.0	Not Detected
Di-n-Butylphthalate	5.0	1.0 J
Fluorene	1.0	Not Detected
Pyrene	1.0	Not Detected
Butylbenzylphthalate	5.0	Not Detected
3,3'-Dichlorobenzidine	20	Not Detected
Chrysene	1.0	Not Detected
Benzo(a)anthracene	1.0	Not Detected
bis(2-Ethylhexyl)phthalate	5.0	Not Detected
Di-n-Octylphthalate	5.0	Not Detected
Benzo(<i>o</i>)fluoranthene	1.0	Not Detected
Benzo(<i>x</i>)fluoranthene	1.0	Not Detected
Benzo(a)pyrene	1.0	Not Detected
Indeno(1,2,3-c,d)pyrene	1.0	Not Detected
Dibenz(a,h)anthracene	1.0	Not Detected
Benzo(g,h)perylene	1.0	Not Detected

J = Estimated value.

Container Type: XAD Tube

Surrogates	%Recovery	Method Limits
2-Fluorophenol	69	50-150
Phenol-d5	72	50-150
Nitrobenzene-d5	75	50-150
2,4,6-Tribromophenol	57	50-150
Fluorene-c10	76	60-120
Pyrene-d10	78	60-120



AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 5 TOX 1 EFF

Lab ID#: 0610414B-05A

MODIFIED EPA METHOD TO-13A GC/MS FULL SCAN

File Name	P102316	Date of Collection	10/19/06
DL Factor	1.00	Date of Analysis	10/23/06 03:17 PM
		Date of Extraction	10/20/06

Compound	Rpt. Limit (ug)	Amount (ug)
Phenol	5.0	Not Detected
bis(2-Chloroethyl) Ether	1.0	Not Detected
4-Chlorophenol	5.0	Not Detected
1,3-Dichlorobenzene	1.0	Not Detected
1,4-Dichlorobenzene	1.0	Not Detected
1,2-Dichlorobenzene	1.0	Not Detected
2-Methylphenol (o-Cresol)	5.0	Not Detected
N-Nitroso-di-n-propylamine	1.0	Not Detected
4-Vinylphenol/3-Vinylphenol	5.0	Not Detected
Hexachloroethane	1.0	Not Detected
Nitrobenzene	1.0	Not Detected
Isophorone	1.0	Not Detected
2-Nitrophenol	5.0	Not Detected
2,4-Dimethylphenol	5.0	Not Detected
bis(2-Chloroethoxy) Methane	1.0	Not Detected
2,4-Dichlorophenol	5.0	Not Detected
1,2,4-Trichlorobenzene	1.0	Not Detected
Naphthalene	1.0	Not Detected
4-Chloroaniline	10	Not Detected
Hexachlorobutadiene	1.0	Not Detected
4-Chloro-3-methylphenol	5.0	Not Detected
2-Methylnaphthalene	1.0	Not Detected
Hexachlorocyclopentadiene	20	Not Detected
2,4,6-Trichlorophenol	5.0	Not Detected
2,4,5-Trichlorophenol	5.0	Not Detected
2-Chloronaphthalene	1.0	Not Detected
2-Nitroaniline	10	Not Detected
Dimethylphthalate	5.0	Not Detected
Acenaphthylene	1.0	Not Detected
2,5-Dinitrotoluene	5.0	Not Detected
3-Nitroaniline	10	Not Detected
Acenaphthene	1.0	Not Detected
2,4-Dinitrophenol	20	Not Detected
4-Nitrophenol	20	Not Detected
2,4-Dinitrotoluene	5.0	Not Detected
Dibenzofuran	1.0	Not Detected
Diethylphthalate	5.0	0.93 J
Fuorene	1.0	Not Detected



AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 5 TOX 1 EFF

Lab ID#: 0610414B-05A

MODIFIED EPA METHOD TO-13A GC/MS FULL SCAN

File Name:	P102315	Date of Collection:	10/19/06
DL Factor:	100	Date of Analysis:	10/23/06 03:17 PM
		Date of Extraction:	10/20/06

Compound	Rpt. Limit (μ g)	Amount (μ g)
4-Chlorophenyl-phenyl Ether	1.0	Not Detected
4-Nitroaniline	10	Not Detected
4,6-Dinitro-2-methylphenol	10	Not Detected
N-Nitrosodiphenylamine	10	Not Detected
4-Bromophenyl-phenyl Ether	1.0	Not Detected
Hexachlorobenzene	1.0	Not Detected
Pentachlorophenol	20	Not Detected
Phenanthrene	1.0	Not Detected
Anthracene	1.0	Not Detected
di-n-Butylphthalate	5.0	Not Detected
Fluoranthene	1.0	Not Detected
Pyrene	1.0	Not Detected
Butylbenzylphthalate	5.0	Not Detected
3,3'-Dichlorobenzidine	20	Not Detected
Chrysene	1.0	Not Detected
Benzo(a)anthracene	1.0	Not Detected
bis(2-Ethylhexyl)phthalate	5.0	Not Detected
Di-n-Octylphthalate	5.0	Not Detected
Benzo(b)fluoranthene	1.0	Not Detected
Benzo(k)fluoranthene	1.0	Not Detected
Benzo(a)pyrene	1.0	Not Detected
Inceno(1,2,3-c,d)pyrene	1.0	Not Detected
Dibenzo(a,h)anthracene	1.0	Not Detected
Benzo(g,h,i)perylene	1.0	Not Detected

.1: Estimated value.

Container Type: XAD Tube

Surrogates	%Recovery	Method Limits
4-Fluorophenol	59	50-150
Phenol-d5	62	50-150
Nitrobenzene-d5	60	50-150
2,4,6-Tribromophenol	56	50-150
Fluorene-d10	61	60-120
Pyrene-d10	68	60-120

OAS
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AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 6 TOX 2 INF

Lab ID#: 0610414B-06A

MODIFIED EPA METHOD TO-13A GC/MS FULL SCAN

File Name:	6TOX2INF	Date of Collection:	10/19/06
DL Factor:	1.00	Date of Analysis:	10/23/06 03:46 PM
			Date of Extraction:
			10/20/06

Compound	Rpt. Limit (ug)	Amount (ug)
Phenol	5.0	Not Detected
bis(2-Chloroethyl) Ether	1.0	3.4
o-Chlorophenol	5.0	Not Detected
1,3-Dichlorobenzene	1.0	1.4
1,4-Dichlorobenzene	1.0	5.1
1,2-Dichlorobenzene	1.0	42
o-Methylphenol (o-Cresol)	5.0	Not Detected
N-Nitroso-di-n-propylamine	1.0	Not Detected
o-Methylphenol/3-Methylphenol	5.0	Not Detected
Hexachloroethane	1.0	Not Detected
Nitrobenzene	1.0	Not Detected
Iso-phorone	1.0	12
2-Nitrophenol	5.0	Not Detected
2,4-Dimethylphenol	5.0	Not Detected
bis(2-Chloroethoxy) Methane	1.0	Not Detected
2,4-Dichlorophenol	5.0	Not Detected
1,2,4-Trichlorobenzene	1.0	1.2
Naphthalene	1.0	24
4-Chloroaniline	10	Not Detected
Hexachlorobutadiene	1.0	2.1
4-Chloro-3-methylphenol	5.0	Not Detected
2-Methylnaphthalene	1.0	4.8
Hexachlorocyclopentadiene	20	Not Detected
2,4,6-Trichlorophenol	5.0	Not Detected
2,4,5-Trichlorophenol	5.0	Not Detected
2-Chloronaphthalene	1.0	Not Detected
2-Nitroaniline	10	Not Detected
O-methylphthalate	5.0	Not Detected
Arenaphthylene	1.0	Not Detected
2,6-Dinitrotoluene	5.0	Not Detected
3-Nitroaniline	10	Not Detected
Arenaphthene	1.0	Not Detected
2,4-Dinitrophenol	20	Not Detected
4-Nitrophenol	20	Not Detected
2,4-Dinitrotoluene	5.0	Not Detected
Dibenzofuran	1.0	Not Detected
Dethylphthalate	5.0	0.81 J 15
Fuorene	1.0	Not Detected

09/17/06



AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 6 TOX 2 INF

Lab ID#: 0610414B-06A

MODIFIED EPA METHOD TO-13A GC/MS FULL SCAN

File Number:	D102317	Date of Collection:	10/19/06
Dil Factor:	1.00	Date of Analysis:	10/23/06 03:48 PM
			Date of Entry:
			10/23/06

Compound	Rpt. Limit (μ g)	Amount (μ g)
4-Chlorophenyl-phenyl Ether	1.0	Not Detected
4-Nitroaniline	10	Not Detected
4,6-Dinitro-2-methylphenol	10	Not Detected
N-N-trosod phenylamine	10	Not Detected
1-Bromophenyl-phenyl Ether	1.0	Not Detected
Hexachlorobenzene	1.0	Not Detected
Pentachlorophenol	20	Not Detected
Phenanthrene	1.0	Not Detected
Anthracene	1.0	Not Detected
Di-n-Butyl phthalate	5.0	Not Detected
Fluoranthene	1.0	Not Detected
Pyrene	1.0	Not Detected
Butylbenzylphthalate	5.0	Not Detected
3,3'-Dichlorobenzidine	20	Not Detected
Chrysene	1.0	Not Detected
Benzo(a)anthracene	1.0	Not Detected
bis(2-Ethylhexyl)phthalate	5.0	12
Di-n-Octylphthalate	5.0	Not Detected
Benzo(b)fluoranthene	1.0	Not Detected
Benzo(k)fluoranthene	1.0	Not Detected
Benzo(a)pyrene	1.0	Not Detected
Indeno(1,2,3-c,d)pyrene	1.0	Not Detected
Dibenz(a,h)anthracene	1.0	Not Detected
benzo(g,h,i)perylene	1.0	Not Detected

J = Estimated value.

Container Type: XAD Tube

Surrogates	%Recovery	Method Limits
2-Fluoropheno!	75	50-150
Phenol-c5	82	50-150
N-tobenzene-d5	79	50-150
2,4,6-Tri bromophenol	64	50-150
Fluorene-d10	74	60-120
Pyrene-c10	77	60-120

OMG
11/16/06



AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 7 TOX 2 INF DUP

Lab ID#: 0610414B-07A

MODIFIED EPA METHOD TO-13A GC/MS FULL SCAN

File Name:	D102318	Date of Collection:	10/19/06
CFR Factor:	1.00	Date of Analysis:	10/23/06 04:16 PM
			Report Date:
			10/20/06

Compound	Rpt. Limit (ug)	Amount (ug)
Phenol	5.0	Not Detected
cis(2-Chloroethyl) Ether	1.0	4.4
2-Chlorophenol	5.0	Not Detected
1,3-Dichlorobenzene	1.0	1.7
1,4-Dichlorobenzene	1.0	6.3
1,2-Dichlorobenzene	1.0	52
2-Methylphenol (c-Cresol)	5.0	Not Detected
N-Nitroso-c-n-propylamine	1.0	Not Detected
4-Methylphenol/3-Methylphenol	5.0	Not Detected
Hexachloroethane	1.0	Not Detected
Nitrobenzene	1.0	Not Detected
Isophorone	1.0	16
2-Naphthol	5.0	Not Detected
2,4-Dimethylphenol	5.0	Not Detected
bis(2-Chloroethoxy) Methane	1.0	Not Detected
2,4-Dichloropheno	5.0	Not Detected
1,2,4-Trichlorobenzene	1.0	1.7
Naphthalene	1.0	32
4-Chloroaniline	10	Not Detected
Hexachlorobutadiene	1.0	2.9
4-Chloro-3-methylphenol	5.0	Not Detected
2-Methylnaphthalene	1.0	6.0
Hexachlorocyclopentadiene	20	Not Detected
2,4,6-Trichloropheno	5.0	Not Detected
2,4,5-Trichloropheno	5.0	Not Detectec
2-Chloronaphthalene	1.0	Not Detected
2-Nitroaniline	10	Not Detected
Diethylphthalate	5.0	Not Detected
Acenaphthylene	1.0	Not Detected
2,6-Dinitrotoluene	5.0	Not Detected
3-Nitroaniline	10	Not Detected
Acenaphthene	1.0	Not Detected
2,4-D nitrophenol	20	Not Detected
4 Nitrophenol	20	Not Detected
2,4-D nitrotoluene	5.0	Not Detected
Dibenzofuran	1.0	Not Detected
Diethylphthalate	5.0	1.1 J
Fuorene	1.0	Not Detected



AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 7 TOX 2 INF DUP

Lab ID#: 0610414B-07A

MODIFIED EPA METHOD TO-13A GC/MS FULL SCAN

File Name:	p102318	Date of Collection:	10/19/06
Oil Factor:	1.00	Date of Analysis:	10/23/06 04:16 PM
		Date of Extractions:	10/20/06

Compound	Rpt. Limit (ug)	Amount (ug)
4-Chlorophenyl-phenyl Ether	1.0	Not Detected
4-Nitroaniline	10	Not Detected
4,6-Dinitro-2-methylphenol	10	Not Detected
N-Nitrosodiphenylamine	10	Not Detected
4-Bromophenyl-phenyl Ether	1.0	Not Detected
Hexachlorobenzene	1.0	Not Detected
Pentachlorophenol	20	Not Detected
Phenanthrene	1.0	Not Detected
Anthracene	1.0	Not Detected
di-n-Butylphthalate	5.0	0.89 J
Fluoranthene	1.0	Not Detected
Pyrene	1.0	Not Detected
Butylbenzylphthalate	5.0	Not Detected
3,3'-Dichlorobenzidine	20	Not Detected
Crylsene	1.0	Not Detected
Benzo(a)anthracene	1.0	Not Detected
bis(2-Ethylhexyl)phthalate	5.0	2.6 J
D-n-Octylphthalate	5.0	Not Detected
Benzo(b)fluoranthene	1.0	Not Detected
Benzo(k)fluoranthene	1.0	Not Detected
Benzo(a)pyrene	1.0	Not Detected
Indeno(1,2,3-c,d)pyrene	1.0	Not Detected
Oibenz(a,b)anthracene	1.0	Not Detected
Benzo(g,h,i)perylene	1.0	Not Detected

J = Estimated value.

Container Type: XAD Tube

Surrogates	%Recovery	Method Limits
2-Fluorophenol	59	50-150
Phenol-d5	82	50-150
4-trobenzene-d5	79	50-150
2,4,6-Tr.bromophenol	65	50-150
Fluorene-d10	75	60-120
Pyrene-d10	77	60-120

10/17/06



AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 8 TOX 2 EFF

Lab ID#: 0610414B-08A

MODIFIED EPA METHOD TO-13A GC/MS FULL SCAN

File Name:	D:\02319\0610414B-08A.DAT	Date of Collection:	10/19/06	
DIL Factor:	1.00	Date of Analysis:	10/23/06 04:46 PM	
			Date of Extraction:	10/20/06

Compound	Rpt. Limit (ug)	Amount (ug)
Phenol	5.0	Not Detected
t-butyl(2-Chloroethyl) Ether	1.0	Not Detected
2-Chlorophenol	5.0	Not Detected
1,3-Dichlorobenzene	1.0	Not Detected
1,4-Dichlorobenzene	1.0	Not Detected
1,2-Dichlorobenzene	1.0	Not Detected
2-Methylphenol (o-Cresol)	5.0	Not Detected
N-Nitroso-di-n-propylamine	1.0	Not Detected
4-Methylphenol/3-Methylphenol	5.0	Not Detected
Hexachloroethane	1.0	Not Detected
Nitrobenzene	1.0	Not Detected
Isochlorone	1.0	Not Detected
2-Nitrophenol	5.0	Not Detected
2,4-Dimethylphenol	5.0	Not Detected
t-butyl(2-Chloroethoxy) Methane	1.0	Not Detected
2,4-Dichlorophenol	5.0	Not Detected
1,2,4-Trichlorobenzene	1.0	Not Detected
Naphthalene	1.0	Not Detected
4-Chloroaniline	10	Not Detected
Hexachlorobutadiene	1.0	Not Detected
4-Chloro-3-methylphenol	5.0	Not Detected
2-Methylnaphthalene	1.0	Not Detected
Hexachlorocyclopentadiene	20	Not Detected
2,4,6-Trichlorophenol	5.0	Not Detected
2,4,5-Trichlorophenol	5.0	Not Detected
2-Chlororaphthalene	1.0	Not Detected
2-Nitroaniline	10	Not Detected
Dimethylphthalate	5.0	Not Detected
Aceanaphthylene	1.0	Not Detected
2,6-Dinitrotoluene	5.0	Not Detected
2-Nitroaniline	10	Not Detected
Aceanaphthene	1.0	Not Detected
2,4-Dinitrophenol	20	Not Detected
4-Nitrophenol	20	Not Detected
2,4-Dinitrotoluene	5.0	Not Detected
Dibenzofuran	1.0	Not Detected
Diethylphthalate	5.0	0.92 J
Fluorene	1.0	Not Detected



AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 8 TOX 2 EFF

Lab ID#: 0610414B-08A

MODIFIED EPA METHOD TO-13A GC/MS FULL SCAN

File Number:	R102319	Date of Collection:	10/19/06
Dil. Factor:	1.00	Date of Analysis:	10/23/06 04:46 PM
			Retention Date of Extraction:
			10/20/06

Compound	Rpt. Limit (ug)	Amount (ug)
4-Chlorophenyl-phenyl Ether	1.0	Not Detected
4-Nitroaniline	10	Not Detected
4,6-Dinitro-2-methylphenol	10	Not Detected
N-Nitrosodiphenylamine	10	Not Detected
4-Bromophenyl-phenyl Ether	1.0	Not Detected
Hexachlorobenzene	1.0	Not Detected
Pentachlorophenol	20	Not Detected
Phenanthrene	1.0	Not Detected
Anthracene	1.0	Not Detected
d-n-Butylphthalate	5.0	Not Detected
Fluoranthene	1.0	Not Detected
Pyrene	1.0	Not Detected
Butylbenzylphthalate	5.0	Not Detected
3,3'-Dichlorobenzidine	20	Not Detected
Chrysene	1.0	Not Detected
Benzo(a)anthracene	1.0	Not Detected
bis(2-Ethylhexyl)phthalate	5.0	1.4 J
D-n-Octylphthalate	5.0	Not Detected
Benzo(b)fluoranthene	1.0	Not Detected
Benzo(k)fluoranthene	1.0	Not Detected
Benzo(a)pyrene	1.0	Not Detected
Indeno(1,2,3-c,d)pyrene	1.0	Not Detected
Dibenz(a,h)anthracene	1.0	Not Detected
Benzo(g,h,i)perylene	1.0	Not Detected

J = Estimated value.

Container Type: XAD Tube

Surrogates	%Recovery	Method Limits
2-Fluorophenol	61	50-150
Phenol-d5	63	50-150
Nitrobenzene-d5	58	50-150
2,4,6-Tr bromophenol	63	50-150
Fluorene-d10	62	60-120
Pyrene-d10	66	60-120

OML
11/17/06

November 2, 2006 Off-Gas Sample Laboratory Results



AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 1 OFFSITE ISVE

Lab ID#: 0611096AR1-01A

MODIFIED EPA METHOD TO-14A GC/MS FULL SCAN

File Name:	111410	Date of Collection:	11/2/06	
Off Factor:	1060	Date of Analysis:	11/14/06 02:58 PM	
Compound	Rpt. Lirnit (ppbv)	Amount (ppbv)	Rpt. Limit (uG/m3)	Amount (uG/m3)
Vinyl Chloride	530	Not Detected	1400	Not Detected
Bromomethane	530	Not Detected	2000	Not Detected
Chloroethane	530	Not Detected	1400	Not Detected
1,1-Dichloroethene	530	1900	2100	7400
Methylene Chloride	530	22000	1800	76000
1,1-Dichloroethane	530	3100	2100	12000
cis-1,2-Dichloroethene	530	1600	2100	6400
Chloroform	530	2200	2600	11000
1,1,1-Trichloroethane	530	25000	2900	130000
Carbon Tetrachloride	530	Not Detected	3300	Not Detected
Benzene	530	13000	1700	43000
1,2-Dichloroethane	530	840	2100	3400
Trichloroethene	530	17000	2800	91000
1,2-Dichloropropane	530	220 J 15	2400	1000 J
cis-1,3-Dichloropropene	530	Not Detected	2400	Not Detected
Toluene	530	110000	2000	420000
trans-1,3-Dichloropropene	530	Not Detected	2400	Not Detected
1,1,2-Trichloroethane	530	180 J 15	2900	1000 J
Tetrachloroethene	530	22000	3600	150000
Chlorobenzene	530	Not Detected	2400	Not Detected
Ethyl Benzene	530	14000	2300	60000
m,p-Xylene	530	57000	2300	250000
o-Xylene	530	21000	2300	90000
Styrene	530	Not Detected	2200	Not Detected
1,1,2,2-Tetrachloroethane	530	Not Detected	3600	Not Detected
Bromochloromethane	530	Not Detected	3600	Not Detected
Dibromochloromethane	530	Not Detected	4500	Not Detected
Chloroethane	2100	Not Detected	4400	Not Detected
Acetone	2100	13000	5000	30000
Carbon Disulfide	2100	Not Detected	6600	Not Detected
trans-1,2-Dichloroethene	2100	Not Detected	8400	Not Detected
2-Butanone (Methyl Ethyl Ketone)	2100	16000	6200	46000
4-Methyl-2-pentanone	2100	7500	8700	31000
2-Hexanone	2100	Not Detected	8700	Not Detected
Bromoform	2100	Not Detected	22000	Not Detected

J = Estimated value.

Container Type: 6 Liter Summa Canister

CVR
12/19/06



AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 1 OFFSITE ISVE

Lab ID#: 0611096AR1-01A

MODIFIED EPA METHOD TO-14A GC/MS FULL SCAN

File Name:	1114410	Date of Collection:	11/12/06
Dil. Factor:	1060	Date of Analysis:	11/14/06 02:50 PM
Surrogates	%Recovery	Method	Limits
1,2-Dichloroethane-d4	105	70-130	
Toluene-d8	100	70-130	
4-Bromofluorobenzene	103	70-130	

OBG
12/19/06



AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 2 SBPA ISVE

Lab ID#: 0611096AR1-02A

MODIFIED EPA METHOD TO-14A GC/MS FULL SCAN

File Name: Dil. Factor:	TO-14A 152	Date of Collection:	11/2/06	
Compound	Rpt. Lirnit (ppbv)	Amount (ppbv)	Rpt. Limit (uG/m3)	Amount (uG/m3)
Vinyl Chloride	180	1800	450	4600
Bromomethane	180	Not Detected	680	Not Detected
Chloroethane	180	380	460	990
1,1-Dichloroethene	180	1100	700	4400
Methylene Chloride	180	5600	610	20000
1,1-Dichloroethane	180	2300	710	9200
cis-1,2-Dichloroethene	180	14000	700	56000
Chloroform	180	4700	860	23000
1,1,1-Trichloroethane	180	18000	960	97000
Carbon Tetrachloride	180	Not Detected	1100	Not Detected
Benzene	180	3500	560	11000
1,2-Dichloroethane	180	290	710	1200
Trichloroethene	180	14000	940	75000
1,2-Dichloropropane	180	270	810	1300
cis-1,3-Dichloropropene	180	Not Detected	800	Not Detected
Toluene	180	39000	660	150000
trans-1,3-Dichloropropene	180	Not Detected	800	Not Detected
1,1,2-Trichloroethane	180	Not Detected	960	Not Detected
Tetrachloroethene	180	26000	1200	180000
Chlorobenzene	180	Not Detected	810	Not Detected
Ethyl Benzene	180	3300	760	32000
m,p-Xylene	180	29000	760	120000
o-Xylene	180	13000	760	56000
Styrene	180	Not Detected	750	Not Detected
1,1,2,2-Tetrachloroethane	180	Not Detected	1200	Not Detected
Bromodichloromethane	180	Not Detected	1200	Not Detected
Dibromochloromethane	180	Not Detected	1500	Not Detected
Chloromethane	700	Not Detected	1400	Not Detected
Acetone	700	1300	1700	3000
Carbon Disulfide	700	Not Detected	2200	Not Detected
trans-1,2-Dichloroethene	700	110 J 15	2800	440 J
2-Butanone (Methyl Ethyl Ketone)	700	810	2100	2400
4-Methyl-2-pentanone	700	1000	2900	4100
2-Hexanone	700	Not Detected	2900	Not Detected
Bromoform	700	Not Detected	7300	Not Detected

J = Estimated value.

Container Type: 6 Liter Summa Canister

CRS
11/19/06
12



AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 2 SBPA ISVE

Lab ID#: 0611096AR1-02A

MODIFIED EPA METHOD TO-14A GC/MS FULL SCAN

File Name:	111411	Date of Collection:	11/2/06
File Extention:	.mz	Date of Analysis:	11/4/06 03:37 PM
Surrogates		%Recovery	Method Limits
1,2-Dichloroethane-d4		102	70-130
Toluene-d8		100	70-130
4-Bromofluorobenzene		109	70-130



AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 3 TOX 1 INF

Lab ID#: 0611096AR1-03A

MODIFIED EPA METHOD TO-14A GC/MS FULL SCAN

File Name Dil Factor	TO-14A 264		Date of Collection: 11/18/06	Date of Analysis: 11/18/06 04:16 PM
Compound	Rpt. Limit (ppbv)	Amount (ppbv)	Rpt. Limit (uG/m3)	Amount (uG/m3)
Vinyl Chloride	130	1200	340	3000
Bromomethane	130	Not Detected	510	Not Detected
Chloroethane	130	180	350	480
1,1-Dichloroethene	130	1100	520	4500
Methylene Chloride	130	6500	460	23000
1,1-Dichloroethane	130	2100	530	8400
cis-1,2-D chloroethene	130	11000	520	44000
Chloroform	130	4600	640	23000
1,1,1-Trichloroethane	130	18000	720	98000
Carbon Tetrachloride	130	Not Detected	830	Not Detected
Benzene	130	2700	420	8800
1,2-Dichloroethane	130	300	530	1200
Trichloroethene	130	15000	710	79000
1,2-Dichloropropane	130	250	610	1200
cis-1,3-Dichloropropene	130	Not Detected	600	Not Detected
Toluene	130	31000	500	120000
trans-1,3-Dichloropropene	130	Not Detected	600	Not Detected
1,1,2-Trichloroethane	130	51 J 15	720	280 J
Tetrachloroethene	130	26000	900	180000
Chlorobenzene	130	Not Detected	610	Not Detected
Ethyl Benzene	130	5600	570	24000
m p-Xylene	130	23000	570	99000
p-Xylene	130	11000	570	46000
Styrene	130	Not Detected	560	Not Detected
1,1,2,2-Tetrachloroethane	130	Not Detected	910	Not Detected
Bromodichlormethane	130	Not Detected	880	Not Detected
Dibromochlormethane	130	Not Detected	1100	Not Detected
Chloromethane	530	Not Detected	1100	Not Detected
Acetone	530	830	1200	2000
Carbon Disulfide	530	Not Detected	1600	Not Detected
trans-1,2-Dichloroethene	530	95 J 15	2100	380 J
2-Butanone (Methyl Ethyl Ketone)	530	580	1600	1700
4-Methyl-2-pentanone	530	660	2200	2700
2-Hexanone	530	Not Detected	2200	Not Detected
Bromoform	530	Not Detected	5400	Not Detected

J = Estimated value.

Container Type: 6 Liter Summa Canister

C76
12/19/06



AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 3 TOX 1 INF

Lab ID#: 0611096AR1-03A

MODIFIED EPA METHOD TO-14A GC/MS FULL SCAN

File Name:	TOX1INF	Date of Collection:	1/17/06
DI Factor:	264	Date of Analysis:	1/14/06 14:16 PM
<hr/>			
Surrogates	%Recovery	Method	Limits
1,2-Dichloroethane-d4	98	70-130	
Toluene-d8	99	70-130	
4-Bromofluorobenzene	111	70-130	

ATL
1/19/06



AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 3 TOX 1 INF Duplicate

Lab ID#: 0611096AR1-03AA

MODIFIED EPA METHOD TO-14A GC/MS FULL SCAN

File Name:	111412	Date of Collection:	11/14/06	
Dil. Factor:	264	Date of Analysis:	11/14/06 04:51 PM	
Compound	Rpt. Limit (ppbv)	Amount (ppbv)	Rpt. Limit (uG/m3)	Amount (uG/m3)
Vinyl Chloride	130	1200	340	2900
Bromomethane	130	Not Detected	510	Not Detected
Chloroethane	130	160	350	410
1,1-Dichloroethene	130	1300	520	5000
Methylene Chloride	130	6200	460	21000
1,1-Dichloroethane	130	2000	530	8000
cis-1,2-Dichloroethene	130	10000	520	42000
Chloroform	130	4400	640	22000
1,1,1-Trichloroethane	130	17000	720	91000
Carbon Tetrachloride	130	Not Detected	830	Not Detected
Benzene	130	2700	420	8600
1,2-Dichloroethane	130	280	530	1200
Trichloroethene	130	14000	710	76000
1,2-Dichloropropane	130	240	610	1100
cis-1,3-Dichloropropene	130	Not Detected	600	Not Detected
Toluene	130	31000	500	120000
trans-1,3-Dichloropropene	130	Not Detected	600	Not Detected
1,1,2-Trichloroethane	130	43 J 15	720	240 J
Tetrachloroethene	130	25000	900	170000
Chlorobenzene	130	Not Detected	610	Not Detected
Ethyl Benzene	130	5500	570	24000
m,p-Xylene	130	22000	570	96000
o-Xylene	130	10000	570	45000
Styrene	130	Not Detected	560	Not Detected
1,1,2,2-Tetrachloroethane	130	Not Detected	910	Not Detected
Bromodichloromethane	130	Not Detected	880	Not Detected
Dibromochloromethane	130	Not Detected	1100	Not Detected
Chlormethane	530	Not Detected	1100	Not Detected
Acetone	530	780	1200	1900
Carbon Disulfide	530	54 J 15	1600	170 J
trans-1,2-Dichloroethene	530	98 J 15	2100	390 J
2-Butanone (Methyl Ethyl Ketone)	530	560	1600	1600
4-Methyl-2-pentanone	530	640	2200	2600
2-Hexanone	530	Not Detected	2200	Not Detected
Bromform	530	Not Detected	5400	Not Detected

J = Estimated value.

Container Type: 6 Liter Summa Canister

12/19/06
JES



AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 3 TOX 1 INF Duplicate

Lab ID#: 0611096AR1-03AA

MODIFIED EPA METHOD TO-14A GC/MS FULL SCAN

File Name:	111141	Date of Collection:	11/12/06
Dil Factor:	264	Date of Analysis:	11/14/06 04:53 PM
Surrogates	%Recovery	Method Limits	
1,2-Dichloroethane-d4	93	70-130	
Toluene-d8	101	70-130	
4-Bromofluorobenzene	109	70-130	

0125
12/19/06



AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 4 TOX 1 INF DUP

Lab ID#: 0611096AR1-04A

MODIFIED EPA METHOD TO-14A GC/MS FULL SCAN

File Name:	111414	Date of Collection:	11/2/06	
DP Factor:	352	Date of Analysis:	11/14/06 /5:30 PM	
Compound	Rpt. Limit (ppbv)	Amount (ppbv)	Rpt. Limit (uG/m3)	Amount (uG/m3)
Vinyl Chloride	180	1800	450	4600
Bromomethane	180	Not Detected	680	Not Detected
Chloroethane	180	500	460	1300
1,1-Dichloroethene	180	840	700	3300
Methylene Chloride	180	5600	610	19000
1,1-Dichloroethane	180	2200	710	8800
cis-1,2-Dichloroethene	180	13000	700	52000
Chloroform	180	4300	860	21000
1,1,1-Trichloroethane	180	16000	960	90000
Carbon Tetrachloride	180	Not Detected	1100	Not Detected
Benzene	180	3200	560	10000
1,2-Dichloroethane	180	260	710	1000
Trichloroethene	180	13000	940	72000
1,2-Dichloropropane	180	250	810	1200
cis-1,3-Dichloropropene	180	Not Detected	800	Not Detected
Toluene	180	33000	660	120000
trans-1,3-Dichloropropene	180	Not Detected	800	Not Detected
1,1,2-Trichloroethane	180	56 J	960	310 J
Tetrachloroethene	180	24000	1200	160000
Chlorobenzene	180	Not Detected	810	Not Detected
Ethyl Benzene	180	6000	760	26000
m,p-Xylene	180	24000	760	100000
o-Xylene	180	10000	760	46000
Styrene	180	Not Detected	750	Not Detected
1,1,2,2-Tetrachloroethane	180	Not Detected	1200	Not Detected
Bromodichloromethane	180	Not Detected	1200	Not Detected
Dibromochloromethane	180	Not Detected	1500	Not Detected
Chloromethane	700	Not Detected	1400	Not Detected
Acetone	700	910	1700	2200
Carbon Disulfide	700	Not Detected	2200	Not Detected
trans-1,2-Dichloroethene	700	100 J	2800	420 J
2-Butanone (Methyl Ethyl Ketone)	700	530 J	2100	1600 J
4-Methyl-2-pentanone	700	880	2900	3600
2-Hexanone	700	Not Detected	2900	Not Detected
Bromoform	700	Not Detected	7300	Not Detected

J = Estimated value.

Container Type: 6 Liter Summa Canister

*OB
12/14/06*



AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 4 TOX 1 INF DUP

Lab ID#: 0611096AR1-04A

MODIFIED EPA METHOD TO-14A GC/MS FULL SCAN

File Name:	111414	Date of Collection:	11/2/06
Lab Folder:	351	Date of Analysis:	11/14/06 02:30 PM

Surrogates	%Recovery	Method Limits
1,2-Dichloroethane-d4	95	70-130
Toluene-d8	101	70-130
4-Bromo fluorobenzene	106	70-130



AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 5 TOX 1 EFF

Lab ID#: 0611096AR1-05A

MODIFIED EPA METHOD TO-14A GC/MS FULL SCAN

File Name:	111415	Date of Collection:	11/2/06	
Eff. Factor:	2.18	Date of Analysis:	11/14/06 06:17 PM	
Compound	Rpt. Limit (ppbv)	Amount (ppbv)	Rpt. Limit (uG/m3)	Amount (uG/m3)
Vinyl Chloride	1.1	91	2.8	230
Bromomethane	1.1	Not Detected	4.2	Not Detected
Chloroethane	1.1	1.6	2.9	4.2
1,1-Dichloroethene	1.1	210	4.3	820
Methylene Chloride	1.1	50	3.8	170
1,1-Dichloroethane	1.1	1.5	4.4	6.0
cis-1,2-Dichloroethene	1.1	98	4.3	390
Chloroform	1.1	7.4	5.3	36
1,1,1-Trichloroethane	1.1	3.4	5.9	18
Carbon Tetrachloride	1.1	1.1	6.8	7.0
Benzene	1.1	170	3.5	560
1,2-Dichloroethane	1.1	0.81 J 15	4.4	3.3 J
Trichloroethene	1.1	160	5.8	890
1,2-Dichloropropane	1.1	Not Detected	5.0	Not Detected
cis-1,3-Dichloropropene	1.1	0.81 J 15	4.9	3.7 J
Toluene	1.1	35	4.1	130
trans-1,3-Dichloropropene	1.1	0.72 J 15	4.9	3.2 J
1,1,2-Trichloroethane	1.1	Not Detected	5.9	Not Detected
Tetrachloroethene	1.1	390	7.4	2600
Chlorobenzene	1.1	8.8	5.0	41
Ethyl Benzene	1.1	6.0	4.7	26
m,p-Xylene	1.1	30	4.7	130
c-Xylene	1.1	11	4.7	50
Styrene	1.1	25	4.6	110
1,1,2,2-Tetrachloroethane	1.1	Not Detected	7.5	Not Detected
Bromodichloromethane	1.1	0.36 J 15	7.3	2.4 J
D bromochloromethane	1.1	Not Detected	9.3	Not Detected
Chloromethane	4.4	15	9.0	32
Acetone	4.4	24	10	57
Carbon Disulfide	4.4	5.4	14	17
trans-1,2-Dichloroethene	4.4	61	17	240
2-Butanone (Methyl Ethyl Ketone)	4.4	8.0	13	24
4-Methyl-2-pentanone	4.4	5.2	18	21
2-Hexanone	4.4	Not Detected	18	Not Detected
Bromoform	4.4	Not Detected	45	Not Detected

J = Estimated value.

Container Type: 6 Liter Summa Canister

10/19/06



AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 5 TOX 1 EFF

Lab ID#: 0611096AR1-05A

MODIFIED EPA METHOD TO-14A GC/MS FULL SCAN

File Name	111415	Date of Collection	11/2/06
DL Factor	2.18	Date of Analysis	11/14/06 06:07 PM

Surrogates	%Recovery	Method Limits
1,2-Dichloroethane-d4	91	70-130
Toluene-d8	100	70-130
4-Bromofluorobenzene	105	70-130



AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 6 TOX 2 INF

Lab ID#: 0611096AR1-06A

MODIFIED EPA METHOD TO-14A GC/MS FULL SCAN

Site Name:	147416		Date of Collection:	1/12/06
DP Factor:	726		Date of Analysis:	1/14/06 06:43 PM
Compound	Rpt. Limit (ppbv)	Amount (ppbv)	Rpt. Limit (uG/m3)	Amount (uG/m3)
Vinyl Chloride	360	550	930	1400
Bromomethane	360	Not Detected	1400	Not Detected
Chloroethane	360	Not Detected	960	Not Detected
1,1-Dichloroethene	360	820	1400	3300
Vinylene Chloride	360	15000	1300	52000
1,1-Dichloroethane	360	2800	1500	11000
cis-1,2-Dichloroethene	360	4900	1400	19000
Chloroforn	360	1700	1800	8400
1,1,1-Trichloroethane	360	20000	2000	110000
Carbon Tetrachloride	360	Not Detected	2300	Not Detected
Benzene	360	11000	1200	36000
1,2-Dichloroethane	360	650	1500	2600
Trichloroethene	360	13000	2000	68000
1,2-Dichloropropane	360	160 J	1700	740 J
cis-1,3-Dichloropropene	360	Not Detected	1600	Not Detected
Toluene	360	85000	1400	320000
trans-1,3-Dichloropropene	360	Not Detected	1600	Not Detected
1,1,2-Trichloroethane	360	130 J	2000	720 J
Tetrachloroethene	360	16000	2500	110000
Chlorobenzene	360	Not Detected	1700	Not Detected
Ethyl Benzene	360	9200	1600	40000
m,p-Xylene	360	38000	1600	170000
o-Xylene	360	14000	1600	59000
Styrene	360	Not Detected	1500	Not Detected
1,1,2,2-Tetrachloroethane	360	Not Detected	2500	Not Detected
Bromodichloromethane	360	Not Detected	2400	Not Detected
Dibromochloromethane	360	Not Detected	3100	Not Detected
Chloromethane	1400	Not Detected	3000	Not Detected
Acetone	1400	15000	3400	36000
Carbon Disulfide	1400	Not Detected	4500	Not Detected
trans-1,2-Dichloroethene	1400	Not Detected	5800	Not Detected
2-Butanone (Methyl Ethyl Ketone)	1400	16000	4300	46000
4-Methyl-2-pentanone	1400	6000	5900	25000
2-Hexanone	1400	Not Detected	5900	Not Detected
Bromoform	1400	Not Detected	15000	Not Detected

J = Estimated value.

Container Type: 6 Liter Summa Canister

CRS
17/1/06



AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 6 TOX 2 INF

Lab ID#: 0611096AR1-06A

MODIFIED EPA METHOD TO-14A GC/MS FULL SCAN

File Name:	File ID:	Date of Collection:	Date of Analysis:
Dil. Factor:			
Surrogates:	%Recovery	Method Limits	
1,2-Dichloroethane-d4	93		70-130
Toluene-d8	100		70-130
4-Bromofluorobenzene	105		70-130

CRS
12/29/06



AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 7 TOX 2 INF DUP

Lab ID#: 0611096AR1-07A

MODIFIED EPA METHOD TO-14A GC/MS FULL SCAN

File Name	111411	Date of Collection	11/2/06	
DP Factor	754	Date of Analysis	11/14/06 07:26 PM	
Compound	Rpt. Limit (ppbv)	Amount (ppbv)	Rpt. Limit (uG/m3)	Amount (uG/m3)
Vinyl Chloride	380	520	960	1300
Bromomethane	380	Not Detected	1500	Not Detected
Chloroethane	380	Not Detected	990	Not Detected
1,1-Dichloroethene	380	910	1500	3600
Methylene Chloride	380	15000	1300	53000
1,1-Dichloroethane	380	2800	1500	11000
cis-1,2-Dichloroethene	380	4700	1500	19000
Chloroform	380	1700	1800	8300
1,1,1-Trichloroethane	380	20000	2000	110000
Carbon Tetrachloride	380	Not Detected	2400	Not Detected
Benzene	380	12000	1200	37000
1,2-Dichloroethane	380	630	1500	2500
Trichloroethene	380	14000	2000	74000
1,2-Dichloropropane	380	210 J	1700	960 J
cis-1,3-Dichloropropene	380	Not Detected	1700	Not Detected
Toluene	380	92000	1400	350000
trans-1,3-Dichloropropene	380	Not Detected	1700	Not Detected
1,1,2-Trichloroethane	380	140 J	2000	770 J
Tetrachloroethene	380	17000	2600	120000
Chlorobenzene	380	Not Detected	1700	Not Detected
Ethyl Benzene	380	10000	1600	45000
m,p-Xylene	380	44000	1600	190000
c-Xylene	380	16000	1600	68000
Styrene	380	Not Detected	1600	Not Detected
1,1,2,2-Tetrachloroethane	380	Not Detected	2600	Not Detected
Bromodichromethane	380	Not Detected	2500	Not Detected
Dibromochromomethane	380	Not Detected	3200	Not Detected
Chloromethane	1500	Not Detected	3100	Not Detected
Acetone	1500	10000	3600	24000
Carbon Disulfide	1500	Not Detected	4700	Not Detected
trans-1,2-Dichloroethene	1500	Not Detected	6000	Not Detected
2-Butanone (Methyl Ethyl Ketone)	1500	13000	4400	38000
2-Methyl-2-pentanone	1500	5800	6200	24000
2-Hexanone	1500	Not Detected	6200	Not Detected
Bromoform	1500	Not Detected	16000	Not Detected

J = Estimated value.

Container Type: 6 Liter Summa Canister

01/06
12/19/06



AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 7 TOX 2 INF DUP

Lab ID#: 0611096AR1-07A

MODIFIED EPA METHOD TO-14A GC/MS FULL SCAN

File Name:	111477	Date of Collection:	11/2/06
TIC Factor:	754	Date of Analysis:	11/14/06 07:28 PM

Surrogates	%Recovery	Method Limits
1,2-Dichloroethane-d4	91	70-130
Toluene-d8	100	70-130
4-Bromo Fluorobenzene	106	70-130

CRS
12/19/06



AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 8 TOX 2 EFF

Lab ID#: 0611096AR1-08A

MODIFIED EPA METHOD TO-14A GC/MS FULL SCAN

File Name Disc Factor	111410 1.3	Date of Collection: 11/21/06	Date of Analysis: 11/14/06 06:04 PM	
Compound	Rpt. Limit (ppbv)	Amount (ppbv)	Rpt. Limit (uG/m3)	Amount (uG/m3)
Vinyl Chloride	6.7	54	17	140
Bromomethane	6.7	Not Detected	26	Not Detected
Chloroethane	6.7	5.6 J 15	18	15 J
1,1-Dichloroethene	6.7	180	26	700
Methylene Chloride	6.7	430	23	1500
1,1-Dichloroethane	6.7	62	27	250
cis-1,2-D chloroethene	6.7	140	26	550
Chloroform	6.7	46	33	220
1,1,1-Trichloroethane	6.7	460	36	2500
Carbon Tetrachloride	6.7	4 1 J 15	42	26 J
Benzene	6.7	490	21	1600
1,2-Dichloroethane	6.7	16	27	65
Trichloroethene	6.7	380	36	2000
1,2-Dichloropropane	6.7	4.1 J 15	31	19 J
cis-1,3-Dichloropropene	6.7	Not Detected	30	Not Detected
Toluene	6.7	1300	25	6900
trans-1,3-Dichloropropene	6.7	Not Detected	30	Not Detected
1,1,2-Trichloroethane	6.7	3.5 J 15	36	19 J
Tetrachloroethene	6.7	640	45	4300
Chlorobenzene	6.7	4.3 J 15	31	20 J
Ethyl Benzene	6.7	180	29	800
m,p-Xylene	6.7	700	29	3000
o-Xylene	6.7	270	29	1200
Styrene	6.7	57	28	240
1,1,2,2-Tetrachloroethane	6.7	Not Detected	46	Not Detected
Bromodichloromethane	6.7	Not Detected	45	Not Detected
Dibromochloromethane	6.7	Not Detected	57	Not Detected
Chloromethane	27	17 J 15	55	35 J
Acetone	27	1100	64	2700
Carbon Disulfide	27	Not Detected	83	Not Detected
trans-1,2-Dichloroethene	27	16 J 15	110	63 J
2-Butanone (Methyl Ethyl Ketone)	27	990	79	2900
4-Methyl-2-pentancne	27	210	110	870
2-Hexanone	27	8.2 J 15	110	33 J
Bromoform	27	Not Detected	280	Not Detected

J = Estimated value.

Container Type: 6 Liter Summa Canister

CNS
12/19/06



AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 8 TOX 2 EFF

Lab ID#: 0611096AR1-08A

MODIFIED EPA METHOD TO-14A GC/MS FULL SCAN

File Name:	111418	Date of Collection:	11/2/06
DP Factor:	13.4	Date of Analysis:	11/19/06 06:04 PM
Surrogates		%Recovery	Method Limits
1,2-Dichloroethane-d4		93	70-130
Toluene-d8		99	70-130
4-Bromofluorobenzene		103	70-130

CPL
12/19/06



AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 1 OFFSITE ISVE

Lab ID#: 0611096B-01A

MODIFIED EPA METHOD TO-13A GC/MS FULL SCAN

File Name:	0111319	Date of Collection:	11/2/06
Div Factor:	1.00	Date of Analysis:	11/13/06 07:28 PM
		Date of Extraction:	11/7/06
Compound	Rpt. Limit (μ g)	Amount (μ g)	
Phenol	5.0		Not Detected
bis(2-Chloroethyl) Ether	1.0		Not Detected
2-Chlorophenol	5.0		Not Detected
1,3-Dichlorobenzene	1.0	2.7	
1,4-Dichlorobenzene	1.0	10	
1,2-Dichlorobenzene	1.0	77	
2-Methylphenol (o-Cresol)	5.0		Not Detected
N-Nitroso-di-n-propylamine	1.0		Not Detected
4-Methylphenol/3-Methylphenol	5.0		Not Detected
Hexachloroethane	1.0		Not Detected
Nitrobenzene	1.0		Not Detected
Isophorone	1.0	41	
2-Nitrophenol	5.0		Not Detected
2,4-Dimethylphenol	5.0		Not Detected
bis(2-Chloroethoxy) Methane	1.0		Not Detected
2,4-Dichlorophenol	5.0		Not Detected
1,2,4-Trichlorobenzene	1.0	4.4	
Naphthalene	1.0	98	
4-Chloraniline	10		Not Detected
Hexachlorobutadiene	1.0	6.4	
4-Chloro-3-methylphenol	5.0		Not Detected
2-Methylnaphthalene	1.0	20	
Hexachlorocyclopentadiene	20		Not Detected
2,4,6-Trichlorophenol	5.0		Not Detected
2,4,5-Trichlorophenol	5.0		Not Detected
2-Chloronaphthalene	1.0		Not Detected
2-Nitroaniline	10		Not Detected
Dimethyl phthalate	5.0		Not Detected
Aceanaphthylene	1.0		Not Detected
2,6-Dinitrotoluene	5.0		Not Detected
3-Nitroaniline	10		Not Detected
Aceanaphthene	1.0		Not Detected
2,4-Dinitrophenol	20		Not Detected
4-Nitrophenol	20		Not Detected
2,4-Dinitrotoluene	5.0		Not Detected
Dibenzofuran	1.0		Not Detected
Diethylphthalate	5.0		Not Detected
Fluorene	1.0		Not Detected



AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 1 OFFSITE ISVE

Lab ID#: 0611096B-01A

MODIFIED EPA METHOD TO-13A GC/MS FULL SCAN

File Name:	0611096B-01A	Date of Collection:	11/2/06
File Ratio:	1.00	Date of Analysis:	11/13/06 12:28 PM
		Date of Extraction:	11/7/06

Compound	Rpt. Limit (μ g)	Amount (μ g)
4-Chlorophenyl-phenyl Ether	1.0	Not Detected
4-Nitroaniline	10	Not Detected
4,6-Dinitro-2-methylphenol	10	Not Detected
N-Nitroso-diphenylamine	10	Not Detected
4-Bromophenyl-phenyl Ether	1.0	Not Detected
Hexachlorobenzene	1.0	Not Detected
Pentachlorophenol	20	Not Detected
Phenanthrene	1.0	Not Detected
Anthracene	1.0	Not Detected
di-n-Butylphthalate	5.0	2.9 J
Fluoranthene	1.0	Not Detected
Pyrene	1.0	Not Detected
Butyl/benzylphthalate	5.0	Not Detected
3,3'-Dichlorobenzidine	20	Not Detected
Chrysene	1.0	Not Detected
Benz(a)anthracene	1.0	Not Detected
bis(2-Ethylhexyl)phthalate	5.0	5.8
Di-n-Octylphthalate	5.0	Not Detected
Benz(b)fluoranthene	1.0	Not Detected
Benz(k)fluoranthene	1.0	Not Detected
Benz(a)pyrene	1.0	Not Detected
Indeno(1,2,3-c,d)pyrene	1.0	Not Detected
Dibenz(a,h)anthracene	1.0	Not Detected
Benz(g,h,i)perylene	1.0	Not Detected

J = Estimated value

Container Type: XAD Tube

Surrogates	%Recovery	Method Limits
2-Fluorophenol	76	50-150
Phenol-d5	78	50-150
Nitrobenzene-d5	87	50-150
2,4,6-Tribromophenol	67	50-150
Fluorene-d10	78	60-120
Pyrene-d10	82	60-120

ONS
12/19/06



AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 2 SBPA ISVE

Lab ID#: 0611096B-02A

MODIFIED EPA METHOD TO-13A GC/MS FULL SCAN

File Name:	0611096B-02A	Date of Collection:	11/09/06
Oil Factor:	1.00	Date of Analysis:	11/10/06 17:58 PM
		Date of Extraction:	11/07/06

Compound	Rpt. Limit (ug)	Amount (ug)
Phenol	5.0	Not Detected
bis(2-Chloroethyl) Ether	1.0	3.2
2-Chlorophenol	5.0	Not Detected
1,3-Dichlorobenzene	1.0	2.7
1,4-Dichlorobenzene	1.0	5.7
1,2-Dichlorobenzene	1.0	24
2-Methylphenol (o-Cresol)	5.0	Not Detected
N-Nitroso-di-n-propylamine	1.0	Not Detected
4-Methylphenol/3-Methylphenol	5.0	Not Detected
Hexachloroethane	1.0	Not Detected
Nitrobenzene	1.0	Not Detected
Isophorone	1.0	2.8
2-Nitrophenol	5.0	Not Detected
2,4-Dimethylphenol	5.0	Not Detected
bis(2-Chloroethoxy) Methane	1.0	Not Detected
2,4-Dichlorophenol	5.0	Not Detected
1,2,4-Trichlorobenzene	1.0	11
Naphthalene	1.0	17
4-Chloroaniline	10	Not Detected
Hexachlorobutadiene	1.0	10
4-Chloro-3-methylphenol	5.0	Not Detected
2-Methylnaphthalene	1.0	9.1
Hexachlorocyclopentadiene	20	Not Detected
2,4,6-Trichlorophenol	5.0	Not Detected
2,4,5-Trichlorophenol	5.0	Not Detected
2-Chloronaphthalene	1.0	Not Detected
2-Nitroaniline	10	Not Detected
Dimethylphthalate	5.0	Not Detected
Acenaphthylene	1.0	Not Detected
2,6-Dinitrotoluene	5.0	Not Detected
3-Nitroaniline	10	Not Detected
Acenaphthene	1.0	Not Detected
2,4-Dinitrophenol	20	Not Detected
4-Nitrophenol	20	Not Detected
2,4-Dinitrotoluene	5.0	Not Detected
Dibenzofuran	1.0	Not Detected
Diethylphthalate	5.0	Not Detected
Fluorene	1.0	Not Detected



AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 2 SBPA ISVE

Lab ID#: 0611096B-02A

MODIFIED EPA METHOD TO-13A GC/MS FULL SCAN

File Name:	D:\EPA\TO-13A\0611096B-02A.DAT	Date of Collection: 11/7/06
File Version:	1.00	Date of Analysis: 11/13/06 U7-SP14M
		Date of Extraction: 11/7/06

Compound	Rpt. Limit (ug)	Amount (ug)
4-Chlorophenyl-phenyl Ether	1.0	Not Detected
4-Nitroaniline	10	Not Detected
4,6-Dinitro-2-methylphenol	10	Not Detected
N-Nitrosodiphenylamine	10	Not Detected
4-Bromophenyl-phenyl Ether	1.0	Not Detected
Hexachlorobenzene	1.0	Not Detected
Pentachlorophenol	20	Not Detected
Phenanthrene	1.0	Not Detected
Anthracene	1.0	Not Detected
di-n-Butylphthalate	5.0	1.4 J 15
Fluoranthene	1.0	Not Detected
Pyrene	1.0	Not Detected
Butylbenzylphthalate	5.0	Not Detected
3,3'-Dichlorobenzidine	20	Not Detected
Chrysene	1.0	Not Detected
Benzo(a)anthracene	1.0	Not Detected
bis(2-Ethylhexyl)phthalate	5.0	4.4 J 15
Di-n-Octylphthalate	5.0	Not Detected
Benzo(b)fluoranthene	1.0	Not Detected
Benzo(k)fluoranthene	1.0	Not Detected
Benzo(a)pyrene	1.0	Not Detected
Indeno(1,2,3-c,d)pyrene	1.0	Not Detected
Dibenz(a,h)anthracene	1.0	Not Detected
Benzoc(g,h,i)perylene	1.0	Not Detected

J = Estimated value.

Container Type: XAD Tube

Surrogates	%Recovery	Method Limits
2-Fluorophenol	80	50-150
Phenol-d5	77	50-150
Nitrobenzene-d5	79	50-150
2,4,6-Tribromophenol	64	50-150
Fluorene-d10	82	60-120
Pyrene-d10	84	60-120



AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 4 TOX1 INF DUP

Lab ID#: 0611096B-03A

MODIFIED EPA METHOD TO-13A GC/MS FULL SCAN

File Name:	4TOX1INF.DAT	Date of Collection:	11/2/06
Off Factor:	1.00	Date of Analysis:	11/13/06 08:26 PM
		Date of Extraction:	11/7/06

Compound	Rpt. Limit (μ g)	Amount (μ g)
Phenol	5.0	Not Detected
bis(2-Choroethyl) Ether	1.0	2.9
2-Chlorophenol	5.0	Not Detected
1,3-Dichlorobenzene	1.0	3.1
1,4-Dichlorobenzene	1.0	6.6
1,2-Dichlorobenzene	1.0	26
2-Methylphenol (α -Cresol)	5.0	Not Detected
N-Nitroso-di-n-propylamine	1.0	Not Detected
4-Methylphenol/3-Methylphenol	5.0	Not Detected
Hexachloroethane	1.0	Not Detected
Nitrobenzene	1.0	Not Detected
Isophorone	1.0	3.0
2-Nitrophenol	5.0	Not Detected
2,4-Dimethylphenol	5.0	Not Detected
bis(2-Chloroethoxy) Methane	1.0	Not Detected
2,4-Dichlorophenol	5.0	Not Detected
1,2,4-Trichlorobenzene	1.0	12
Naphthalene	1.0	18
4-Chlcroaniline	10	Not Detected
Hexachlorobutadiene	1.0	11
4-Chlro-3-methylphenol	5.0	Not Detected
2-Methylnaphthalene	1.0	9.6
Hexachlorocyclopentadiene	20	Not Detected
2,4,6-Trichlorophenol	5.0	Not Detected
2,4,5-Trichlorophenol	5.0	Not Detected
2-Chlronaphthalene	1.0	Not Detected
2-Nitroaniline	1.0	Not Detected
Dimethylphthalate	5.0	Not Detected
Acenaphthylene	1.0	Not Detected
2,6-Dinitrotoluene	5.0	Not Detected
3-Nitroaniline	1.0	Not Detected
Acenaphthene	1.0	Not Detected
2,4-Dinitrophenol	20	Not Detected
4-Nitrophenol	20	Not Detected
2,4-Dinitrotoluene	5.0	Not Detected
Dibenzofuran	1.0	Not Detected
Diethyl phthalate	5.0	Not Detected
Fluorene	1.0	Not Detected

OAS
12/19/06



AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 4 TOX1 INF DUP

Lab ID#: 0611096B-03A

MODIFIED EPA METHOD TO-13A GC/MS FULL SCAN

File Name:	0411321	Date of Collection:	11/27/06
Dil. Factor:	1.00	Date of Analysis:	11/13/06 08:28 PM
		Date of Extraction:	11/7/06

Compound	Rpt. Limit (μ g)	Amount (μ g)
4-Chlorophenyl-phenyl Ether	1.0	Not Detected
4-Nitroaniline	10	Not Detected
4,6-Dinitro-2-methylphenol	10	Not Detected
N-Nitrosodiphenylamine	10	Not Detected
4-Bromophenyl-phenyl Ether	1.0	Not Detected
Hexachlorobenzene	1.0	Not Detected
Pentachlorophenol	20	Not Detected
Phenanthrene	1.0	Not Detected
Anthracene	1.0	Not Detected
di-n-Butylphthalate	5.0	1.1 J /5
Fluoranthene	1.0	Not Detected
Pyrene	1.0	Not Detected
Butylbenzylphthalate	5.0	Not Detected
3,3'-Dichlorobenzidine	20	Not Detected
Chrysene	1.0	Not Detected
Benzo(a)anthracene	1.0	Not Detected
bis(2-Ethylhexyl)phthalate	5.0	1.5 J /5
Di-n-Octylphthalate	5.0	Not Detected
Benzo(b)fluoranthene	1.0	Not Detected
Benzo(k)fluoranthene	1.0	Not Detected
Benzo(a)pyrene	1.0	Not Detected
Indeno(1,2,3-c,d)pyrene	1.0	Not Detected
Dibenz(a,h)anthracene	1.0	Not Detected
Benzo(g,h,i)perylene	1.0	Not Detected

J = Estimated value.

Container Type: XAD Tube

Surrogates	%Recovery	Method Limits
2-Fluorophenol	78	50-150
Phenol-d5	74	50-150
Nitrobenzene-d5	80	50-150
2,4,6-Tribromophenol	59	50-150
Fluorene-d10	80	60-120
Pyrene-d10	80	60-120

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12/19/06



AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 5 TOX1 EFF

Lab ID#: 0611096B-04A

MODIFIED EPA METHOD TO-13A GC/MS FULL SCAN

File Number:	P110322	Date of Collection: 4/12/06
Oil Factor:	1.00	Date of Analysis: 4/17/06 08:56 PM
		Date of Extraction: 4/17/06

Compound	Rpt. Limit (ug)	Amount (ug)
Phenol	5.0	Not Detected
bis(2-Chloroethyl) Ether	1.0	Not Detected
2-Chlorophenol	5.0	Not Detected
1,3-Dichlorobenzene	1.0	Not Detected
1,4-Dichlorobenzene	1.0	Not Detected
1,2-Dichlorobenzene	1.0	Not Detected
2-Methylphenol (o-Cresol)	5.0	Not Detected
N-Nitroso-di-n-propylamine	1.0	Not Detected
4-Methylphenol/3-Methylphenol	5.0	Not Detected
Hexachloroethane	1.0	Not Detected
Nitrobenzene	1.0	Not Detected
Isophorone	1.0	Not Detected
2-Nitrophenol	5.0	Not Detected
2,4-D-methylphenol	5.0	Not Detected
bis(2-Chloroethoxy) Methane	1.0	Not Detected
2,4-Dichlorophenol	5.0	Not Detected
1,2,4-Trichlorobenzene	1.0	Not Detected
Naphthalene	1.0	Not Detected
4-Chloraniline	10	Not Detected
Hexachlorobutadiene	1.0	Not Detected
4-Chloro-3-methylphenol	5.0	Not Detected
2-Methylnaphthalene	1.0	Not Detected
Hexachlorocyclopentadiene	20	Not Detected
2,4,6-Trichlorophenol	5.0	Not Detected
2,4,5-Trichlorophenol	5.0	Not Detected
2-Chloronaphthalene	1.0	Not Detected
2-Nitroaniline	10	Not Detected
Dimethylphthalate	5.0	Not Detected
Acenaphthylene	1.0	Not Detected
2,6-Dinitrotoluene	5.0	Not Detected
3-Nitroaniline	10	Not Detected
Acenaphthene	1.0	Not Detected
2,4-Dinitrophenal	20	Not Detected
4-Nitrophenol	20	Not Detected
2,4-Dinitrotoluene	5.0	Not Detected
Dibenzofuran	1.0	Not Detected
Diethylphthalate	5.0	Not Detected
Fluorene	1.0	Not Detected



AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 5 TOX1 EFF

Lab ID#: 0611096B-04A

MODIFIED EPA METHOD TO-13A GC/MS FULL SCAN

File Name:	5TOX1EFF.DAT	Date of Collection:	11/2/06
Dil Factor:	1.00	Date of Analysis:	11/13/06 08:51 PM
		Date of Extraction:	11/2/06

Compound	Rpt. Limit (μ g)	Amount (μ g)
4-Chlorophenyl-phenyl Ether	1.0	Not Detected
4-Nitroaniline	10	Not Detected
4,6-Dinitro-2-methylphenol	10	Not Detected
N-Nitrosodiphenylamine	10	Not Detected
4-Bromophenyl-phenyl Ether	1.0	Not Detected
Hexachlorobenzene	1.0	Not Detected
Pentachlorophenol	20	Not Detected
Phenanthrene	1.0	Not Detected
Anthracene	1.0	Not Detected
di-n-Butylphthalate	5.0	1.3 J
Fluoranthene	1.0	Not Detected
Pyrene	1.0	Not Detected
Butylbenzylphthalate	5.0	Not Detected
3,3'-D chlorobenzidine	20	Not Detected
Chrysene	1.0	Not Detected
Benz(a)anthracene	1.0	Not Detected
bis(2-Ethylhexyl)phthalate	5.0	Not Detected
Di-n-Octylphthalate	5.0	Not Detected
Benz(b)fluoranthene	1.0	Not Detected
Benz(k)fluoranthene	1.0	Not Detected
Benz(a)pyrene	1.0	Not Detected
Indeno(1,2,3-c,d)pyrene	1.0	Not Detected
Dibenz(a,h)anthracene	1.0	Not Detected
Benz(g,h,i)perylene	1.0	Not Detected

J = Estimated value.

Container Type: XAD Tube

Surrogates	%Recovery	Method Limits
2-Fluorophenol	73	50-150
Phenol-d5	74	50-150
Nitrobenzene-d5	73	50-150
2,4,6-Tribromophenol	63	50-150
Fluorene-d10	72	60-120
Pyrene-d10	77	60-120

ONS
12/19/06



AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 7 TOX 2 INF DUP

Lab ID#: 0611096B-05A

MODIFIED EPA METHOD TO-13A GC/MS FULL SCAN

File Name:	0611096B-05A	Date of Collection:	11/2/06
Dil Factor:	1.00	Date of Analysis:	11/13/06 09:24 PM
		Date of Extraction:	11/7/06

Compound	Rpt. Limit (μ g)	Amount (μ g)
Phenol	5.0	Not Detected
bis(2-Chloroethyl) Ether	1.0	2.2
2-Chlorophenol	5.0	Not Detected
1,3-Dichlorobenzene	1.0	1.0
1,4-Dichlorobenzene	1.0	3.2
1,2-Dichlorobenzene	1.0	24
2-Methylphenol (o-Cresol)	5.0	Not Detected
N-Nitroso-di-n-propylamine	1.0	Not Detected
4-Methylphenol/3-Methylphenol	5.0	Not Detected
Hexachloroethane	1.0	Not Detected
Nitrobenzene	1.0	Not Detected
Isophorone	1.0	5.3
2-Nitrophenol	5.0	Not Detected
2,4-Dimethylphenol	5.0	Not Detected
bis(2-Chloroethoxy) Methane	1.0	Not Detected
2,4-Dichlorophenol	5.0	Not Detected
1,2,4-Trichlorobenzene	1.0	0.67 J 15
Naphthalene	1.0	12
4-Chloroaniline	10	Not Detected
Hexachlorobutadiene	1.0	1.2
4-Chloro-3-methylphenol	5.0	Not Detected
2-Methylnaphthalene	1.0	2.2
Hexachlorocyclooctadiene	20	Not Detected
2,4,6-Trichlorophenol	5.0	Not Detected
2,4,5-Trichlorophenol	5.0	Not Detected
2-Chloronaphthalene	1.0	Not Detected
2-Nitroaniline	10	Not Detected
Dimethylphthalate	5.0	Not Detected
Acenaphthylene	1.0	Not Detected
2,6-Dinitrotoluene	5.0	Not Detected
3-Nitroaniline	10	Not Detected
Acenaphthene	1.0	Not Detected
2,4-Dinitrophenol	20	Not Detected
4-Nitrophenol	20	Not Detected
2,4-Dinitrotoluene	5.0	Not Detected
Dibenzofuran	1.0	Not Detected
Diethylphthalate	5.0	Not Detected
Fluorene	1.0	Not Detected

CHS
12/19/06



AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 7 TOX 2 INF DUP

Lab ID#: 0611096B-05A

MODIFIED EPA METHOD TO-13A GC/MS FULL SCAN

File Name:	R111323	Date of Collection:	11/2/06
Dil Factor:	1.00	Date of Analysis:	11/7/06 04:27 PM
		Date of Extraction:	11/2/06

Compound	Rpt. Limit (μ g)	Amount (μ g)
4-Chlorophenyl-phenyl Ether	1.0	Not Detected
4-Nitroaniline	10	Not Detected
4,6-Dinitro-2-methylphenol	10	Not Detected
N-Nitrosodiphenylamine	10	Not Detected
4-Bromophenyl-phenyl Ether	1.0	Not Detected
Hexachlorobenzene	1.0	Not Detected
Pentachlorophenol	20	Not Detected
Phenanthrene	1.0	Not Detected
Anthracene	1.0	Not Detected
di-n-Butylphthalate	5.0	1.1 J 15
Fluoranthene	1.0	Not Detected
Pyrene	1.0	Not Detected
Butylbenzylphthalate	5.0	Not Detected
3,3'-Dichlorobenzidine	20	Not Detected
Chrysene	1.0	Not Detected
Benz(a)anthracene	1.0	Not Detected
bis(2-Ethylhexyl)phthalate	5.0	1.1 J 15
Di-n-Octylphthalate	5.0	Not Detected
Benzo(b)fluoranthene	1.0	Not Detected
Benzo(k)fluoranthene	1.0	Not Detected
Benzo(a)pyrene	1.0	Not Detected
Indeno(1,2,3-c,d)pyrene	1.0	Not Detected
Dibenz(a,h)anthracene	1.0	Not Detected
Benz(g,h,i)perylene	1.0	Not Detected

J = Estimated value.

Container Type: XAD Tube

Surrogates	%Recovery	Method Limits
2-Fluorophenol	80	50-150
Phenol-d5	77	50-150
Nitrobenzene-d5	85	50-150
2,4,6-Tribromophenol	61	50-150
Fluorene-d10	79	60-120
Fyrene-d10	82	60-120

*APC
12/11/06*



AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 8 TOX 2 EFF

Lab ID#: 0611096B-06A

MODIFIED EPA METHOD TO-13A GC/MS FULL SCAN

File Name:	0111124	Date of Collection:	11/2/06
Dil Factor:	1.00	Date of Analysis:	11/3/06 09:57 PM
		Date of Extraction:	11/7/06

Compound	Rpt. Limit (ug)	Amount (ug)
Phenol	5.0	Not Detected
bis(2-Chloroethyl) Ether	1.0	Not Detected
2-Chlorophenol	5.0	Not Detected
1,3-Dichlorobenzene	1.0	Not Detected
1,4-Dichlorobenzene	1.0	Not Detected
1,2-Dichlorobenzene	1.0	1.4
2-Methylphenol (o-Cresol)	5.0	Not Detected
N-Nitroso-di-n-propylamine	1.0	Not Detected
4-Methylphenol/3-Methylphenol	5.0	Not Detected
Hexachloroethane	1.0	Not Detected
Nitrobenzene	1.0	Not Detected
Isophorone	1.0	Not Detected
2-Nitrophenol	5.0	Not Detected
2,4-D methylphenol	5.0	Not Detected
bis(2-Chloroethoxy) Methane	1.0	Not Detected
2,4-Dichlorophenol	5.0	Not Detected
1,2,4-Trichlorobenzene	1.0	Not Detected
Naphthalene	1.0	2.0
4-Chloroaniline	10	Not Detected
Hexachlorobutadiene	1.0	Not Detected
4-Chloro-3-methylphenol	5.0	Not Detected
2-Methylnaphthalene	1.0	Not Detected
Hexachlorocyclopentadiene	20	Not Detected
2,4,6-Trichlorophenol	5.0	Not Detected
2,4,5-Trichlorophenol	5.0	Not Detected
2-Chloronaphthalene	1.0	Not Detected
2-Nitroaniline	10	Not Detected
Dimethylphthalate	5.0	Not Detected
Aceanaphthylene	1.0	Not Detected
2,6-Dinitrotoluene	5.0	Not Detected
3-Nitroaniline	10	Not Detected
Aceranaphthene	1.0	Not Detected
2,4-Dinitrophenol	20	Not Detected
4-Nitrophenol	20	Not Detected
2,4-Dinitrotoluene	5.0	Not Detected
Dibenzofuran	1.0	Not Detected
Diethylphthalate	5.0	Not Detected
Fluorene	1.0	Not Detected



AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 8 TOX 2 EFF

Lab ID#: 0611096B-06A

MODIFIED EPA METHOD TO-13A GC/MS FULL SCAN

File Name:	1111324	Date of Collection:	11/7/06
Dil. Factor:	1.00	Date of Analysis:	11/12/06 09:57 PM
		Date of Extraction:	11/7/06

Compound	Rpt. Limit (ug)	Amount (ug)
4-Chlorophenyl-phenyl Ether	1.0	Not Detected
4-Nitroariline	1.0	Not Detected
4,6-Dinitro-2-methylphenol	1.0	Not Detected
N-Nitrosodiphenylamine	1.0	Not Detected
4-Bromophenyl-phenyl Ether	1.0	Not Detected
Hexachlorobenzene	1.0	Not Detected
Pentachlorophenol	20	Not Detected
Phenanthrene	1.0	Not Detected
Anthracene	1.0	Not Detected
di-n-Butylphthalate	5.0	1.1 J
Fluoranthene	1.0	Not Detected
Pyrene	1.0	Not Detected
Butylbenzylphthalate	5.0	Not Detected
3,3'-D chlorobenzidine	20	Not Detected
Chrysene	1.0	Not Detected
Benz(a)anthracene	1.0	Not Detected
bis(2-Ethylhexyl)phthalate	5.0	3.8 J
Di-n-Cetylphthalate	5.0	Not Detected
Benz(b)fluoranthene	1.0	Not Detected
Benz(k)fluoranthene	1.0	Not Detected
Benz(a)pyrene	1.0	Not Detected
Indeno(1,2,3-c,d)pyrene	1.0	Not Detected
Dibenz(a,h)anthracene	1.0	Not Detected
Benzo(c,h,i)perylene	1.0	Not Detected

J = Estimated value.

Container Type: XAD Tube

Surrogates	%Recovery	Method Limits
2-Fluorophenol	67	50-150
Phenol-d5	70	50-150
Nitrobenzene-d5	70	50-150
2,4,6-Tribromophenol	53	50-150
Fluorophenol-d10	73	60-120
Pyrene-d10	75	60-120

CRS
12/19/06

December 11, 2006 Off-Gas Sample Laboratory Results



AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 1 OFFSITE ISVE

Lab ID#: 0612226AR1-01A

MODIFIED EPA METHOD TO-14A GC/MS FULL SCAN

File Number EN Factor	Sample ID	Date of Collection: 12/17/06 Date of Analysis: 12/25/06 05:05 PM		
Compound	Rpt. Limit (ppbv)	Amount (ppbv)	Rpt. Limit (uG/m3)	Amount (uG/m3)
Vinyl Chloride	270	Not Detected	680	Not Detected
Bromomethane	270	Not Detected	1000	Not Detected
Chloroethane	270	Not Detected	710	Not Detected
1,1-Dichloroethene	270	360	1100	1400
Methylene Chloride	270	19000	930	65000
1,1-Dichloroethane	270	2700	1100	11000
cis-1,2-Dichloroethene	270	1400	1100	5600
Chloroform	270	1800	1300	8700
1,1,1-Trichloroethane	270	21000	1500	110000
Carbon Tetrachloride	270	Not Detected	1700	Not Detected
Benzene	270	12000	860	38000
1,2-Dichloroethane	270	660	1100	2700
Trichloroethene	270	13000	1400	69000
1,2-Dichloropropane	270	180 J 15	1200	830 J
cis-1,3-Dichloropropene	270	Not Detected /C	1200	Not Detected
Toluene	270	86000	1000	320000
trans-1,3-Dichloropropene	270	Not Detected	1200	Not Detected
1,1,2-Trichloroethane	270	160 J 15	1500	880 J
Tetrachloroethene	270	16000	1800	110000
Chlorobenzene	270	Not Detected	1200	Not Detected
Ethyl Benzene	270	10000	1200	45000
m,p-Xylene	270	44000	1200	190000
c-Xylene	270	17000	1200	72000
Styrene	270	Not Detected	1100	Not Detected
1,1,2,2-Tetrachloroethane	270	Not Detected	1800	Not Detected
Bromodichloromethane	270	Not Detected	1800	Not Detected
Dibromochloromethane	270	Not Detected	2300	Not Detected
Chloromethane	1100	Not Detected	2200	Not Detected
Acetone	1100	13000	2500	39000
Carbon Disulfide	1100	440 J 15	3300	1400 J
trans-1,2-Dichloroethene	1100	Not Detected	4200	Not Detected
2-Butanone (Methyl Ethyl Ketone)	1100	16000	3200	49000
4-Methyl-2-pentanone	1100	7100	4400	29000
2-Hexanone	1100	220 J 15	4400	890 J
Bromoform	1100	Not Detected	11000	Not Detected

J = Estimated value.

Container Type: 6 Liter Summa Canister

CH
2/6/07



AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 1 OFFSITE ISVE

Lab ID#: 0612226AR1-01A

MODIFIED EPA METHOD TO-14A GC/MS FULL SCAN

Surrogates	%Recovery	Method Limits
1,2-Dichloroethane-d4	96	70-130
Toluene-d8	100	70-130
4-Bromofluorobenzene	97	70-130



AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 2 SBPA ISVE

Lab ID#: 0612226AR1-02A

MODIFIED EPA METHOD TO-14A GC/MS FULL SCAN

Compound	Rpt. Limit (ppbv)	Amount (ppbv)	Rpt. Limit (uG/m3)	Amount (uG/m3)
Vinyl Chloride	130	1300	340	3300
Bromomethane	130	Not Detected	520	Not Detected
Chloroethane	130	640 J	350	1700
1,1-Dichloroethene	130	1100	530	4200
Methylene Chloride	130	6400	460	22000
1,1-Dichloroethane	130	2900	540	12000
cis-1,2-Dichloroethene	130	14000	530	56000
Chloroform	130	4600	650	22000
1,1,1-Trichloroethane	130	17000	730	95000
Carbon Tetrachloride	130	Not Detected	840	Not Detected
Benzene	130	5200	430	17000
1,2-Dichloroethane	130	250	540	1000
Trichloroethene	130	13000	720	72000
1,2-Dichloropropane	130	290	620	1400
cis-1,3-Dichloropropene	130	Not Detected	610	Not Detected
Toluene	130	42000	500	160000
trans-1,3-Dichloropropene	130	Not Detected	610	Not Detected
1,1,2-Trichloroethane	130	66 J J	730	360 J
Tetrachloroethene	130	25000	910	170000
Chlorobenzene	130	57 J J	620	260 J
Ethyl Benzene	130	10000	580	45000
m,p-Xylene	130	44000	580	190000
o-Xylene	130	20000	580	88000
Styrene	130	Not Detected	570	Not Detected
1,1,2,2-Tetrachloroethane	130	Not Detected	920	Not Detected
Bromodichloromethane	130	Not Detected	900	Not Detected
Dibromochloromethane	130	Not Detected	1100	Not Detected
Chloromethane	540	Not Detected	1100	Not Detected
Acetone	540	1200	1300	2800
Carbon Disulfide	540	77 J J	1700	240 J
trans-1,2-Dichloroethene	540	120 J J	2100	470 J
2-Butanone (Methyl Ethyl Ketone)	540	1000	1600	3000
4-Methyl-2-pentanone	540	1600	2200	6400
2-Hexanone	540	Not Detected	2200	Not Detected
Bromoform	540	Not Detected	5500	Not Detected

J = Estimated value.

Container Type: 6 Liter Summa Canister

CH
2/6/07



AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 2 SBPA ISVE

Lab ID#: 0612226AR1-02A

MODIFIED EPA METHOD TO-14A GC/MS FULL SCAN

Surrogate	%Recovery	Method Limits
1,2-Dichloroethane-d4	82	70-130
Toluene-d8	105	70-130
4-Bromofluorobenzene	91	70-130



AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 3 TOX 1 INF

Lab ID#: 0612226AR1-03A

MODIFIED EPA METHOD TO-14A GC/MS FULL SCAN

File Number	11-22-12-264	Date of Collection	11/22/12	
Sample Date	264	Date of Analysis	11/22/12 06:00 PM	
Compound	Rpt. Limit (ppbv)	Amount (ppbv)	Rpt. Limit (uG/m3)	Amount (uG/m3)
Vinyl Chloride	130	1500	340	3800
Bromomethane	130	Not Detected	520	Not Detected
Chloroethane	130	770 <i>J</i>	350	2000
1,1-Dichloroethene	130	1200	530	5000
Methylene Chloride	130	6600	460	23000
1,1-Dichloroethane	130	3100	540	12000
cis-1,2-Dichloroethene	130	15000	530	60000
Chloroform	130	4900	650	24000
1,1,1-Trichloroethane	130	18000	730	100000
Carbon Tetrachloride	130	Not Detected	840	Not Detected
Benzene	130	4800	430	15000
1,2-Dichloroethane	130	270	540	1100
Trichloroethene	130	14000	720	74000
1,2-Dichloropropane	130	300	620	1400
cis-1,3-D chloropropene	130	Not Detected	610	Not Detected
Toluene	130	41000	500	150000
trans-1,3-D chloropropene	130	Not Detected	610	Not Detected
t,1,2-Trichloroethane	130	47 J <i>J</i>	730	260 J
Tetrachloroethene	130	26000	910	170000
Chlorobenzene	130	69 J <i>J</i>	620	320 J
Ethyl Benzene	130	9300	580	40000
m,p-Xylene	130	39000	580	170000
p-Xylene	130	18000	580	77000
Styrene	130	Not Detected	570	Not Detected
1,1,2,2-Tetrachloroethane	130	Not Detected	920	Not Detected
Bromodichloromethane	130	Not Detected	900	Not Detected
Dibromochloromethane	130	Not Detected	1100	Not Detected
Chlorormethane	540	Not Detected	1100	Not Detected
Acetone	540	1200	1300	2900
Carbon Disulfide	540	33 J <i>J</i>	1700	100 J
trans-1,2-Dichloroethene	540	110 J <i>J</i>	2100	440 J
2-Butanone (Methyl Ethyl Ketone)	540	940	1600	2800
1-Methyl-2-pentanone	540	1500	2200	6300
3-Hexanone	540	Not Detected	2200	Not Detected
Bromoform	540	Not Detected	5500	Not Detected

J = Estimated value.

Container Type: 6 Liter Summa Canister

*OK
2/1/13*



AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 3 TOX 1 INF

Lab ID#: 0612226AR1-03A

MODIFIED EPA METHOD TO-14A GC/MS FULL SCAN

File Name:	0612226AR1-03A	Date of Collection:	12/11/08
DIC Factor:	-60	Date of Analysis:	12/22/08 10:11 PM
Surrogates	%Recovery	Method Limits	
1,2-Dichloroethane-d4	81	70-130	
Toluene-d8	104	70-130	
4-Bromofluorobenzene	91	70-130	



AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 4 TOX 1 INF DUP

Lab ID#: 0612226AR1-04A

MODIFIED EPA METHOD TO-14A GC/MS FULL SCAN

File Number Dil. Factor	Date Collected PPB	Date of Analysis PPM	Rpt. Limit (uG/m3)	Amount (uG/m3)
Compound	Rpt. Lirnit (ppbv)	Amount (ppbv)		
Vinyl Chloride	130	1500	340	3900
Bromomethane	130	Not Detected	520	Not Detected
Chloroethane	130	780 J	350	2000
1,1-Dichloroethene	130	1300	530	5200
Methylene Chloride	130	7500	460	26000
1,1-Dichloroethane	130	3300	540	14000
cis-1,2-Dichloroethene	130	17000	530	66000
Chloroform	130	5300	650	26000
1,1,1-Trichloroethane	130	19000	730	100000
Carbon Tetrachloride	130	Not Detected	840	Not Detected
Benzene	130	5600	430	18000
1,2-Dichloroethane	130	280	540	1100
Trichloroethene	130	15000	720	82000
1,2-Dichloropropane	130	340	620	1600
cis-1,3-Dichloropropene	130	Not Detected	610	Not Detected
Toluene	130	43000	500	180000
trans-1,3-Dichloropropene	130	Not Detected	610	Not Detected
1,1,2-Trichloroethane	130	70 J J	730	380 J
Tetrachloroethene	130	28000	910	190000
Chlorobenzene	130	97 J J	620	440 J
Ethyl Benzene	130	11000	580	48000
m,p-Xylene	130	46000	580	200000
o-Xylene	130	21000	580	91000
Styrene	130	Not Detected	570	Not Detected
1,1,2,2-Tetrachloroethane	130	Not Detected	920	Not Detected
Bromodichloromethane	130	Not Detected	900	Not Detected
Dibromochloromethane	130	Not Detected	1100	Not Detected
Chloromethane	540	Not Detected	1100	Not Detected
Acetone	540	1300	1300	3000
Carbon Disulfide	540	39 J J	1700	120 J
trans-1,2-Dichloroethene	540	140 J J	2100	560 J
2-Butanone (Methyl Ethyl Ketone)	540	1100	1600	3200
4-Methyl-2-pentanone	540	1800	2200	7400
2-Hexanone	540	Not Detected	2200	Not Detected
Bromoform	540	Not Detected	5500	Not Detected

J = Estimated value.

Container Type: 6 Liter Summa Canister

*OK
JUL 11/11*



AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 4 TOX 1 INF DUP

Lab ID#: 0612226AR1-04A

MODIFIED EPA METHOD TO-14A GC/MS FULL SCAN

File Number:	M22214	Date of Collection:	12/17/06
Dil. Factor:	266	Date of Analysis:	12/22/06 07:31 PM
Surrogates	%Recovery	Method Limits	
1,2-Dichloroethane-d4	79	70-130	
Toluene-d8	105	70-130	
4-Bromofluorobenzene	89	70-130	

CMH
JLW



AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 5 TOX 1 EFF

Lab ID#: 0612226AR1-05A

MODIFIED EPA METHOD TO-14A GC/MS FULL SCAN

File Name	Sample ID	Date of Sample	Report Date	
Oil Factor	1.00	1/22/2016 10:12 AM	1/22/2016 10:12 PM	
Compound	Rpt. Limit (ppbv)	Amount (ppbv)	Rpt. Limit (uG/m3)	Amount (uG/m3)
Vinyl Chloride	0.67	76	1.7	190
Bromomethane	0.67	Not Detected	2.6	Not Detected
Chloroethane	0.67	3.2 J 15	1.8	8.4
1,1-Dichloroethene	0.67	200	2.6	800
Methylene Chloride	0.67	50	2.3	170
1,1-Dichloroethane	0.67	7.9	2.7	32
cis-1,2-Dichloroethene	0.67	87	2.6	340
Chloroform	0.67	13	3.3	62
1,1,1-Trichloroethane	0.67	43	3.6	240
Carbon Tetrachloride	0.67	0.60 J 15	4.2	3.8 J
Benzene	0.67	120	2.1	400
1,2-Dichloroethane	0.67	0.85	2.7	3.4
Trichloroethene	0.67	120	3.6	630
1,2-Dichloropropane	0.67	0.77	3.1	3.6
cis-1,3-D chloropropene	0.67	0.25 J 15	3.0	1.1 J
Toluene	0.67	100	2.5	390
trans-1,3-Dichloropropene	0.67	0.24 J 15	3.0	1.1 J
1,1,2-Trichloroethane	0.67	0.25 J 15	3.6	1.4 J
Tetrachloroethene	0.67	260	4.5	1800
Chlorobenzene	0.67	6.3	3.1	29
Ethyl Benzene	0.67	24	2.9	110
m,p-Xylene	0.67	110	2.9	480
o-Xylene	0.67	50	2.9	220
Slyrene	0.67	12	2.8	51
1,1,2,2-Tetrachloroethane	0.67	Not Detected	4.6	Not Detected
Bromodichloromethane	0.67	Not Detected	4.5	Not Detected
D bromochloromethane	0.67	Not Detected	5.7	Not Detected
Chloromethane	2.7	8.9	5.5	18
Acetone	2.7	23	6.4	55
Carbon Disulfide	2.7	120	8.3	380
trans-1,2-Dichloroethene	2.7	43	11	170
2-Butanone (Methyl Ethyl Ketone)	2.7	560 E 1E	7.9	1600 E
4-Methyl-2-pentanone	2.7	44	11	180
2-Hexanone	2.7	1.2 J 15	11	4.7 J
Bromoform	2.7	Not Detected	28	Not Detected

J = Estimated value.

E = Exceeds instrument calibration range.

Container Type: 6 Liter Summa Canister

CHS
2/26/07



AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 5 TOX 1 EFF

Lab ID#: 0612226AR1-05A

MODIFIED EPA METHOD TO-14A GC/MS FULL SCAN

File Name:	0612226AR1-05A	Date of Collection:	12/22/06
RI Factor:	0.94	Initial Analysis:	12/22/06 09:12 PM

Surrogates	%Recovery	Method Limits
1,2-Dichloroethane-d4	76	70-130
Toluene-d8	106	70-130
4-Bromofluorobenzene	94	70-130

CH₂
2/16/07



AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 6 TOX 2 INF

Lab ID#: 0612226AR1-06A

MODIFIED EPA METHOD TO-14A GC/MS FULL SCAN

File Name or Project	12/22/10 520	Date of Collection	12/22/10 Date of Analysis	12/23/10 10:05:31 PM
Compound	Rpt. Limit (ppbv)	Amount (ppbv)	Rpt. Limit (uG/m3)	Amount (uG/m3)
Vinyl Chloride	260	Not Detected	660	Not Detected
Bromomethane	260	Not Detected	1000	Not Detected
Chloroethane	260	Not Detected	690	Not Detected
1,1-Dichloroethene	260	380	1000	1500
Methylene Chloride	260	19000	900	66000
1,1-Dichloroethane	260	2800	1000	11000
cis-1,2-Dichloroethene	260	1400	1000	5600
Chloroform	260	1800	1300	8800
1,1,1-Trichloroethane	260	22000	1400	120000
Carbon Tetrachloride	260	Not Detected	1600	Not Detected
Benzene	260	13000	830	40000
1,2-Dichloroethane	260	650	1000	2600
Trichloroethylene	260	13000	1400	70000
1,2-Dichloropropane	260	200 J 15	1200	920 J
cis-1,3-Dichloropropene	260	Not Detected 18	1200	Not Detected
Toluene	260	83000	980	310000
trans-1,3-Dichloropropene	260	Not Detected	1200	Not Detected
1,1,2-Trichloroethane	260	140 J 15	1400	770 J
Tetrachloroethylene	260	16000	1800	110000
Chlorobezene	260	Not Detected	1200	Not Detected
Ethyl Benzene	260	9700	1100	42000
m,p-Xylene	260	40000	1100	180000
o-Xylene	260	15000	1100	64000
Styrene	260	Not Detected	1100	Not Detected
1,1,2,2-Tetrachloroethane	260	Not Detected	1800	Not Detected
Bromoc dichloromethane	260	Not Detected	1700	Not Detected
Dibromochloromethane	260	Not Detected	2200	Not Detected
Chloromethane	1000	Not Detected	2100	Not Detected
Acetone	1000	15000	2500	35000
Carbon Disulfide	1000	170 J 15	3200	540 J
trans-1,2-Dichloroethene	1000	Not Detected	4100	Not Detected
2-Butanone (Methyl Ethyl Ketone)	1000	16000	3100	46000
4-Methyl-2-pentanone	1000	6600	4300	27000
2-Hexanone	1000	180 J 15	4300	750 J
Bromoform	1000	Not Detected	11000	Not Detected

J = Estimated value.

Container Type: 6 Liter Summa Canister

CH
2/6/11



AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 6 TOX 2 INF

Lab ID#: 0612226AR1-06A

MODIFIED EPA METHOD TO-14A GC/MS FULL SCAN

Surrogates	RT	Date of Collection	Date of Analysis	Method Limits
1,2-Dichloroethane-d4	720	1/17/06	1/23/06 10:57 AM	70-130
Toluene-d8				70-130
4-Bromo Fluorobenzene				70-130



AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 6 TOX 2 INF Duplicate

Lab ID#: 0612226AR1-06AA

MODIFIED EPA METHOD TO-14A GC/MS FULL SCAN

File Name DL Factor	1/25/01 120	Date of Collection: 1/25/01	Date of Analysis: 1/25/01 08:48 PM	
Compound	Rpt. Limit (ppbv)	Amount (ppbv)	Rpt. Limit (uG/m3)	Amount (uG/m3)
Vinyl Chloride	260	Not Detected	660	Not Detected
Bromomethane	260	Not Detected	1000	Not Detected
Chloroethane	260	Not Detected	690	Not Detected
1,1-Dichloroethene	260	380	1000	1500
Methylene Chloride	260	19000	900	66000
1,1-Dichloroethane	260	2800	1000	11000
cis-1,2-Dichloroethene	260	1400	1000	5500
Chloroform	260	1900	1300	9100
1,1,1-Trichloroethane	260	22000	1400	120000
Carbon Tetrachloride	260	Not Detected	1600	Not Detected
Benzene	260	13000	830	41000
1,2-Dichloroethane	260	720	1000	2900
Trichloroethene	260	13000	1400	71000
1,2-Dichloropropane	260	210 J 15	1200	980 J
cis-1,3-Dichloropropene	260	Not Detected AC	1200	Not Detected
Toluene	260	84000	980	320000
trans-1,3-Dichloropropene	260	Not Detected	1200	Not Detected
1,1,2-Trichloroethane	260	150 J 15	1400	830 J
Tetrachloroethene	260	17000	1800	110000
Chlorobenzene	260	Not Detected	1200	Not Detected
Ethyl Benzene	260	10000	1100	43000
m,p-Xylene	260	42000	1100	180000
o-Xylene	260	15000	1100	67000
Styrene	260	Not Detected	1100	Not Detected
1,1,2,2-Tetrachloroethane	260	Not Detected	1800	Not Detected
Bromodichloromethane	260	Not Detected	1700	Not Detected
Dibromochloromethane	260	Not Detected	2200	Not Detected
Chloromethane	1000	Not Detected	2100	Not Detected
Acetone	1000	15000	2500	36000
Carbon Disulfide	1000	150 J 15	3200	470 J
trans-1,2-Dichloroethene	1000	Not Detected	4100	Not Detected
2-Butanone (Methyl Ethyl Ketone)	1000	16000	3100	48000
4-Methyl-2-pentanone	1000	7000	4300	28000
2-Hexanone	1000	180 J 15	4300	720 J
Bromoform	1000	Not Detected	11000	Not Detected

J = Estimated value.

Container Type: 6 Liter Summa Canister

APC
3/6/01



AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 6 TOX 2 INF Duplicate

Lab ID#: 0612226AR1-06AA

MODIFIED EPA METHOD TO-14A GC/MS FULL SCAN

File Number:	162226	Date of Collection:	12/1/06
Dil. Factor:	54.0	Date of Analysis:	12/5/06 06:34 PM

Surrogates	%Recovery	Method Limits
1,2-Dichloroethane-d4	96	70-130
Toluene-d8	99	70-130
4-Bromofluorobenzene	98	70-130



AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 7 TOX 2 INF DUP

Lab ID#: 0612226AR1-07A

MODIFIED EPA METHOD TO-14A GC/MS FULL SCAN

File Number	47225-10 340	Date of Collection	12/17/06	Date of Analysis	12/20/06 10:47 PM
Compound	Rpt. Limit (ppbv)	Amount (ppbv)	Rpt. Limit (uG/m3)	Amount (uG/m3)	
Vinyl Chloride	260	Not Detected	660	Not Detected	
Bromomethane	260	Not Detected	1000	Not Detected	
Chloroethane	260	Not Detected	690	Not Detected	
1,1-Dichloroethene	260	430	1000	1700	
Methylene Chloride	260	18000	900	63000	
1,1-Dichloroethane	260	2700	1000	11000	
cis-1,2-Dichloroethene	260	1300	1000	5100	
Chloroform	260	1800	1300	8600	
1,1,1-Trichloroethane	260	21000	1400	120000	
Carbon Tetrachloride	260	Not Detected	1600	Not Detected	
Benzene	260	12000	830	38000	
1,2-Dichloroethane	260	650	1000	2600	
Trichloroethene	260	12000	1400	67000	
1,2-Dichloropropane	260	190 J	1200	890 J	
cis-1,3-Dichloropropene	260	Not Detected	1200	Not Detected	
Toluene	260	79000	980	300000	
trans-1,3-D chloropropene	260	Not Detected	1200	Not Detected	
1,1,2-Trichloroethane	260	120 J	1400	640 J	
Tetrachloroethene	260	16000	1800	110000	
Chlorobenzene	260	Not Detected	1200	Not Detected	
Ethyl Benzene	260	9100	1100	40000	
m,p-Xylene	260	38000	1100	160000	
c-Xylene	260	14000	1100	59000	
Styrene	260	Not Detected	1100	Not Detected	
1,1,2,2-Tetrachloroethane	260	Not Detected	1800	Not Detected	
Bromodichloromethane	260	Not Detected	1700	Not Detected	
D-bromoethane	260	Not Detected	2200	Not Detected	
Chloromethane	1000	Not Detected	2100	Not Detected	
Acetone	1000	8400	2500	20000	
Carbon Disulfide	1000	140 J	3200	430 J	
trans-1,2-Dichloroethene	1000	Not Detected	4100	Not Detected	
2-Butanone (Methyl Ethyl Ketone)	1000	2000	3100	6000	
4-Methyl-2-pentanone	1000	2700	4300	11000	
2-Hexanone	1000	Not Detected	4300	Not Detected	
Bromoform	1000	Not Detected	11000	Not Detected	

J = Estimated value.

Container Type: 6 Liter Summa Canister

CH
2007



AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 7 TOX 2 INF DUP

Lab ID#: 0612226AR1-07A

MODIFIED EPA METHOD TO-14A GC/MS FULL SCAN

File Name Dil. Ratio	Sample ID Series	Date of Collection Time	Date of Analysis Time	Method Limits
Surrogates		%Recovery		
1,2-Dichloroethane-d4		96		70-130
Toluene-d8		101		70-130
4-Bromofluorobenzene		101		70-130



AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 7 TOX 2 INF DUP Duplicate

Lab ID#: 0612226AR1-07AA

MODIFIED EPA METHOD TO-14A GC/MS FULL SCAN

Compound	Rpt. Limit (ppbv)	Amount (ppbv)	Rpt. Limit (uG/m3)	Amount (uG/m3)
Vinyl Chloride	260	Not Detected	660	Not Detected
Bromomethane	260	Not Detected	1000	Not Detected
Chloroethane	260	Not Detected	690	Not Detected
1,1-Dichloroethene	260	500	1000	2000
Methylene Chloride	260	18000	900	61000
1,1-Dichloroethane	260	2600	1000	11000
cis-1,2-Dichloroethene	260	1200	1000	4900
Chloroform	260	1700	1300	8400
1,1,1-Trichloroethane	260	20000	1400	110000
Carbon Tetrachloride	260	Not Detected	1600	Not Detected
Benzene	260	12000	830	38000
1,2-Dichloroethane	260	600	1000	2400
Trichloroethene	260	12000	1400	66000
1,2-Dichloropropane	260	180 J	1200	810 J
cis-1,3-Dichloropropene	260	Not Detected	1200	Not Detected
Toluene	260	78000	980	290000
trans-1,3-Dichloropropene	260	Not Detected	1200	Not Detected
1,1,2-Trichloroethane	260	120 J	1400	680 J
Tetrachloroethene	260	15000	1800	100000
Chlorobenzene	260	Not Detected	1200	Not Detected
Ethyl Benzene	260	8900	1100	39000
m,p-Xylene	260	37000	1100	160000
n-Xylene	260	13000	1100	58000
Styrene	260	Not Detected	1100	Not Detected
1,1,2,2-Tetrachloroethane	260	Not Detected	1800	Not Detected
Bromochloromethane	260	Not Detected	1700	Not Detected
Dibromochloromethane	260	Not Detected	2200	Not Detected
Chloroform	1000	Not Detected	2100	Not Detected
Acetone	1000	8000	2500	19000
Carbon Disulfide	1000	130 J	3200	420 J
trans-1,2-Dichloroethene	1000	Not Detected	4100	Not Detected
2-Butanone (Methyl Ethyl Ketone)	1000	1700	3100	5100
4-Methyl-2-pentanone	1000	2800	4300	11000
2-Hexanone	1000	Not Detected	4300	Not Detected
Bromoform	1000	Not Detected	11000	Not Detected

J = Estimated value.

Container Type: 6 Liter Summa Canister

CHS
2/6/07



AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 7 TOX 2 INF DUP Duplicate

Lab ID#: 0612226AR1-07AA

MODIFIED EPA METHOD TO-14A GC/MS FULL SCAN

Surrogates	%Recovery	Method Limits
1,2-Dichloroethane-d4	96	70-130
Toluene-d8	100	70-130
4-Bromofluorobenzene	100	70-130



AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 8 TOX 2 EFF

Lab ID#: 0612226AR1-08A

MODIFIED EPA METHOD TO-14A GC/MS FULL SCAN

File Name: DIF Factor:	112226AR1 1.0	Date of Collection: 12/2/2006	Date of Analysis: 12/2/2006 03:57 PM	
Compound	Rpt. Limit (ppbv)	Amount (ppbv)	Rpt. Limit (uG/m3)	Amount (uG/m3)
Vinyl Chloride	9.0	40	23	100
Bromomethane	9.0	Not Detected	35	Not Detected
Chloroethane	9.0	Not Detected	24	Not Detected
1,1-Dichloroethene	9.0	190	35	750
Methylene Chloride	9.0	870	31	3000
1,1-Dichloroethane	9.0	110	36	450
cis-1,2-Dichloroethene	9.0	81	35	320
Chloroform	9.0	77	44	380
1,1,1-Trichloroethane	9.0	720	49	3900
Carbon Tetrachloride	9.0	Not Detected	56	Not Detected
Benzene	9.0	820	28	2600
1,2-Dichloroethane	9.0	25	36	100
Trichloroethene	9.0	550	48	3000
1,2-Dichloropropane	9.0	7.2 J	41	34 J
cis-1,3-Dichloropropene	9.0	Not Detected	41	Not Detected
Toluene	9.0	2900	34	11000
trans-1,3-Dichloropropene	9.0	Not Detected	41	Not Detected
1,1,2-Trichloroethane	9.0	5.4 J	49	29 J
Tetrachloroethene	9.0	780	61	5300
Chlorobenzene	9.0	3.5 J	41	16 J
Ethyl Benzene	9.0	270	39	1200
m,p-Xylene	9.0	1100	39	4700
o-Xylene	9.0	400	39	1700
Styrene	9.0	150	38	650
1,1,2,2-Tetrachloroethane	9.0	2.8 J	61	19 J
Bromodichloromethane	9.0	Not Detected	60	Not Detected
Dibromochloromethane	9.0	Not Detected	76	Not Detected
Chloromethane	36	18 J	74	38 J
Acetone	36	770	85	1800
Carbon Disulfide	36	9.8 J	110	31 J
trans-1,2-Dichloroethene	36	12 J	140	46 J
2-Butanone (Methyl Ethyl Ketone)	36	500	100	1500
2-Methyl-2-pentanone	36	120	150	500
2-Hexanone	36	5.7 J	150	23 J
Bromoform	36	Not Detected	370	Not Detected

J = Estimated value.

Container Type: 6 Liter Summa Canister



AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 8 TOX 2 EFF

Lab ID#: 0612226AR1-08A

MODIFIED EPA METHOD TO-14A GC/MS FULL SCAN

File Number	Sample ID	Date of Collection	Date of Analysis	Method Limits
0612226	8 TOX 2 EFF	12/14/06	12/21/06 01:57 PM	
Surrogates	%Recovery			
1,2-Dichloroethane-d4	88			70-130
Toluene-d8	102			70-130
4-Bromof uorobenzene	90			70-130

01/5
2/1/10



AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: I OFFSITE ISVE

Lab ID#: 0612226B-01A

MODIFIED EPA METHOD TO-13A GC/MS FULL SCAN

File Name:	Sample ID:	Date of Collection:
Alt. Sample ID:	Sample Type:	Date of Analysis:
		Date of Extraction:

Compound	Rpt. Limit (ug)	Amount (ug)
Phenol	5.0	5.2
bis(2-Chloroethyl) Ether	1.0	6.2
2-Chlorophenol	5.0	Not Detected
1,3-Dichlorobenzene	1.0	2.5
1,4-Dichlorobenzene	1.0	7.9
1,2-Dichlorobenzene	1.0	65
2-Methoxyphenol (o-Cresol)	5.0	Not Detected
N-Nitroso-di-n-propylamine	1.0	Not Detected
4-Methoxyphenol/3-Methylphenol	5.0	3.3 J 15
Hexachloroethane	1.0	Not Detected
Nitrobenzene	1.0	Not Detected
Isophorone	1.0	42
2-Nitrophenol	5.0	Not Detected
2,4-D methylphenol	5.0	Not Detected
bis(2-Chloroethoxy) Methane	1.0	Not Detected
2,4-Dichlorophenol	5.0	Not Detected
1,2,4-Trichlorobenzene	1.0	3.6
Naphthalene	1.0	76
4-Chloroaniline	10	Not Detected
Hexachlorobutadiene	1.0	4.8
4-Chloro-3-methylphenol	5.0	Not Detected
2-Methylnaphthalene	1.0	16
Hexachlorocyclopentadiene	20	Not Detected
2,4,6-Trichlorophenol	5.0	Not Detected
2,4,5-Trichlorophenol	5.0	Not Detected
2-Chloronaphthalene	1.0	Not Detected
2-Nitroaniline	10	Not Detected
Dimethylphthalate	5.0	Not Detected
Acenaphthylene	1.0	Not Detected
2,6-Dinitrotoluene	5.0	Not Detected
3-Nitroaniline	10	Not Detected
Acenaphthene	1.0	Not Detected
2,4-Dinitrophenol	20	Not Detected
4-Nitrophenol	20	Not Detected
2,4-Dinitrotoluene	5.0	Not Detected
Dibenzofuran	1.0	Not Detected
Diethylphthalate	5.0	Not Detected
Fluorene	1.0	Not Detected

2/6/07



AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 1 OFFSITE ISVE

Lab ID#: 0612226B-01A

MODIFIED EPA METHOD TO-13A GC/MS FULL SCAN

File Number:	P121410	Date of Collection:	12/14/06
Dil. Factor:	1.00	Date of Analysis:	12/14/06 02:44 PM
		Date of Extraction:	12/14/06

Compound	Rpt. Limit (μ g)	Amount (μ g)
4-Chlorophenyl-phenyl Ether	1.0	Not Detected
4-Nitroaniline	10	Not Detected
4,6-Dinitro-2-methylphenol	10	Not Detected
N-Nitrosodiphenylamine	10	Not Detected
4-Bromophenyl-phenyl Ether	1.0	Not Detected
Hexachlorobenzene	1.0	Not Detected
Pentachlorophenol	20	Not Detected
Phenanthrene	1.0	Not Detected
Anthracene	1.0	Not Detected
d-n-Butylphthalate	5.0	Not Detected
Fluoranthene	1.0	Not Detected
Pyrene	1.0	Not Detected
Butylbenzylphthalate	5.0	Not Detected
3,3'-Dichlorobenzidine	20	Not Detected
Chrysene	1.0	Not Detected
Benzo(a)anthracene	1.0	Not Detected
bis(2-Ethylhexyl)phthalate	5.0	3.5 J
Di-n-Octylphthalate	5.0	Not Detected
Benzo(b)fluoranthene	1.0	Not Detected
Benzo(k)fluoranthene	1.0	Not Detected
Benzo(a)pyrene	1.0	Not Detected
Indeno(1,2,3-c,d)pyrene	1.0	Not Detected
Dibenz(a,h)anthracene	1.0	Not Detected
Benz(c,g,h,i)perylene	1.0	Not Detected

J = Estimated value.

Q = Exceeds Quality Control limits.

Container Type: XAD Tube

Surrogates	%Recovery	Method Limits
2-Fluorophenol	10 Q	50-150
Phenol-d5	100	50-150
Nitrobenzene-d5	84	50-150
2,4,6-Tribromophenol	75	50-150
Fluorene-d10	78	60-120
Pyrene-d10	83	60-120

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2/6/07



AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 1 OFFSITE ISVE Duplicate

Lab ID#: 0612226B-01AA

MODIFIED EPA METHOD TO-13A GC/MS FULL SCAN

File Name:	0612226B-01AA	Date of Collection:	12/14/06
Dil Factor:	1.00	Date of Analysis:	12/14/06 03:14 PM
		Date of Extraction:	12/12/06

Compound	Rpt. Limit (μ g)	Amount (μ g)
Phenol	5.0	5.4
bis(2-Chloroethyl) Ether	1.0	6.7
2-Chlorophenol	5.0	Not Detected
1,3-Dichlorobenzene	1.0	2.6
1,4-Dichlorobenzene	1.0	8.4
1,2-Dichlorobenzene	1.0	69
2-Methylphenol (o-Cresol)	5.0	Not Detected
N-Nitroso-di-n-propylamine	1.0	Not Detected
4-Methylphenol/3-Methylphenol	5.0	3.5 J 15
Hexachloroethane	1.0	Not Detected
Nitrobenzene	1.0	Not Detected
Isophorone	1.0	43
2-Nitrophenol	5.0	Not Detected
2,4-Dimethylphenol	5.0	Not Detected
bis(2-Chloroethoxy) Methane	1.0	Not Detected
2,4-Dichlorophenol	5.0	Not Detected
1,2,4-Trichlorobenzene	1.0	3.4
Naphthalene	1.0	77
4-Chloroaniline	10	Not Detected
Hexachlorobutadiene	1.0	5.1
4-Chloro-3-methylphenol	5.0	Not Detected
2-Methyl naphthalene	1.0	17
Hexachlorocyclopentadiene	20	Not Detected
2,4,6-Tr chlorophenol	5.0	Not Detected
2,4,5-Tr chlorophenol	5.0	Not Detected
2-Chloronaphthalene	1.0	Not Detected
2-Nitroaniline	10	Not Detected
Dimethylphthalate	5.0	Not Detected
Acenaphthylene	1.0	Not Detected
2,6-D nitrotoluene	5.0	Not Detected
3-Nitroaniline	10	Not Detected
Acenaphthene	1.0	Not Detected
2,4-Dinitrophenol	20	Not Detected
4-Nitrophenol	20	Not Detected
2,4-Dinitrotoluene	5.0	Not Detected
Dibenzofuran	1.0	Not Detected
Diethylphthalate	5.0	Not Detected
Fluorene	1.0	Not Detected



AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 1 OFFSITE ISVE Duplicate

Lab ID#: 0612226B-01AA

MODIFIED EPA METHOD TO-13A GC/MS FULL SCAN

File Name:	0612226B-01AA	Date of GC Collection:	12/14/06
Off. Factor:	1.00	Date of Analysis:	12/14/06 04:14 PM
		Date of Extraction:	12/14/06

Compound	Rpt. Limit (ug)	Amount (ug)
4-Chlorophenyl-phenyl Ether	1.0	Not Detected
4-Nitroaniline	10	Not Detected
4,6-Dinitro-2-methylphenol	10	Not Detected
N-Nitrosodiphenylamine	10	Not Detected
4-Bromophenyl-phenyl Ether	1.0	Not Detected
Hexachlorobenzene	1.0	Not Detected
Pentachlorophenol	20	Not Detected
Phenanthrene	1.0	Not Detected
Anthracene	1.0	Not Detected
di-n-Butylphthalate	5.0	Not Detected
Fluoranthene	1.0	Not Detected
Pyrene	1.0	Not Detected
Butylbenzylphthalate	5.0	Not Detected
3,3'-Dichlorobenzidine	20	Not Detected
Chrysene	1.0	Not Detected
Benz[a]anthracene	1.0	Not Detected
b s(2-Ethylhexyl)phthalate	5.0	3.7 J 15
Di-n-Octylphthalate	5.0	Not Detected
Benz[b]fluoranthene	1.0	Not Detected
Benz[k]fluoranthene	1.0	Not Detected
Benz(a)pyrene	1.0	Not Detected
Indenc(1,2,3-c,d)pyrene	1.0	Not Detected
Dibenz(a,h)anthracene	1.0	Not Detected
Benzo(g,h,i)perylene	1.0	Not Detected

J = Estimated value.

Q = Exceeds Quality Control limits.

Container Type: XAD Tube

Surrogates	%Recovery	Method Limits
2-Fluorophenol	10 Q	50-150
Phencl-d5	106	50-150
Nitrobenzene-d5	86	50-150
2,4,6-Tribromophenol	80	50-150
Fluorene-d10	83	60-120
Pyrene-d10	86	60-120

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AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 2 SBPA.ISVE

Lab ID#: 0612226B-02A

MODIFIED EPA METHOD TO-13A GC/MS FULL SCAN

File Number:	DTZ1441	Date of Collection:	12/14/06
DL Factor:	1.00	Date of Analysis:	12/14/06 13:42:PM
		Date of Extraction:	12/14/06

Compound	Rpt. Limit (μ g)	Amount (μ g)
Phenol	5.0	Not Detected
bis(2-Chloroethyl) Ether	1.0	3.2
2-Chlorophenol	5.0	Not Detected
1,3-Dichlorobenzene	1.0	3.4
1,4-Dichlorobenzene	1.0	6.9
1,2-Dichlorobenzene	1.0	28
2-Methylphenol (o-Cresol)	5.0	Not Detected
N-Nitroso-di-n-propylamine	1.0	Not Detected
4-Methylphenol/3-Methylphenol	5.0	Not Detected
Hexachloroethane	1.0	Not Detected
N-trobenzene	1.0	Not Detected
Isophorone	1.0	3.0
2-Nitrophenol	5.0	Not Detected
2,4-Dinethylphenol	5.0	Not Detected
bis(2-Chloroethoxy) Methane	1.0	Not Detected
2,4-Dichlorophenol	5.0	Not Detected
1,2,4-Trichlorobenzene	1.0	19
Naphthalene	1.0	26
4-Chloroaniline	10	Not Detected
Hexachlorobutadiene	1.0	13
4-Chloro-3-methylphenol	5.0	Not Detected
2-Methylnaphthalene	1.0	16
Hexachlorocyclopentadiene	20	Not Detected
2,4,6-Trichlorophenol	5.0	Not Detected
2,4,5-Trichlorophenol	5.0	Not Detected
2-Chlronaphthalene	1.0	Not Detected
2-Nitroaniline	10	Not Detected
Dimethylphthalate	5.0	Not Detected
Acenaphthylene	1.0	Not Detected
2,6-Dinitrotoluene	5.0	Not Detected
3-Nitroaniline	10	Not Detected
Acenaphthene	1.0	Not Detected
2,4-Dinitrophenol	20	Not Detected
4-Nitrophenol	20	Not Detected
2,4-Dinitrotoluene	5.0	Not Detected
Dibenzofuran	1.0	Not Detected
Diethylphthalate	5.0	Not Detected
Fluorene	1.0	Not Detected



AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 2 SBP A ISVE

Lab ID#: 0612226B-02A

MODIFIED EPA METHOD TO-13A GC/MS FULL SCAN

File Name:	G121412	Date of Collection:	12/14/06
File Date:	12/14/06	Date of Analysis:	12/14/06 03:44 PM
		Date of Extraction:	12/12/06

Compound	Rpt. Limit (μ g)	Amount (μ g)
4-Chlorophenyl-phenyl Ether	1.0	Not Detected
4-Nitroaniline	10	Not Detected
4,6-Dinitro-2-methylphenol	10	Not Detected
N-Nitrosodiphenylamine	10	Not Detected
4-Bromophenyl-phenyl Ether	1.0	Not Detected
Hexachlorobenzene	1.0	Not Detected
Pentachlorophenol	20	Not Detected
Phenanthrene	1.0	Not Detected
Anthracene	1.0	Not Detected
di-n-Butylphthalate	5.0	Not Detected
Fluoranthene	1.0	Not Detected
Pyrene	1.0	Not Detected
Butylbenzylphthalate	5.0	Not Detected
3,3'-Dichlorobenzidine	20	Not Detected
Chrysene	1.0	Not Detected
Benz(a)anthracene	1.0	Not Detected
bis(2-Ethylhexyl)phthalate	5.0	1.8 J
Di-n-Octylphthalate	5.0	Not Detected
Benz(b)fluoranthene	1.0	Not Detected
Benz(k)fluoranthene	1.0	Not Detected
Benz(a)pyrene	1.0	Not Detected
Indeno(1,2,3-c,d)pyrene	1.0	Not Detected
Dibenz(a,h)anthracene	1.0	Not Detected
Benzo(ℓ ,h,i)perylene	1.0	Not Detected

J = Estimated value.

Q = Exceeds Quality Control limits.

Container Type: XAD Tube

Surrogates	%Recovery	Method Limits
2-Fluorophenol	39 Q	50-150
Phenol-d5	100	50-150
Nitrobenzene-d5	89	50-150
2,4,6-Tribromophenol	84	50-150
Fluorene-d10	90	60-120
Pyrene-d10	99	60-120

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AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 3 TOX 1 INF

Lab ID#: 0612226B-03A

MODIFIED EPA METHOD TO-13A GC/MS FULL SCAN

File Number:	0612226B-03A	Date of Collection:	12/14/06
File Date:	12/14/06	Date of Analysis:	12/14/06 04:15 PM
		Date of Extraction:	12/14/06

Compound	Rpt. Limit (ug)	Amount (ug)
Phenol	5.0	Not Detected
bis(2-Chloroethyl) Ether	1.0	2.6
2-Chlorophenol	5.0	Not Detected
1,3-Dichlorobenzene	1.0	2.8
1,4-Dichlorobenzene	1.0	5.8
1,2-Dichlorobenzene	1.0	23
2-Methyphenol (o-Cresol)	5.0	Not Detected
N-Nitroso-di-n-propylamine	1.0	Not Detected
4-Methylphenol/3-Methylphenol	5.0	Not Detected
Hexachloroethane	1.0	Not Detected
Nitrobenzene	1.0	Not Detected
Isophorone	1.0	2.4
2-Nitrophenol	5.0	Not Detected
2,4-Dimethylphenol	5.0	Not Detected
bis(2-Chloroethoxy) Methane	1.0	Not Detected
2,4-Dichlorophenol	5.0	Not Detected
1,2,4-Trichlorobenzene	1.0	13
Naphthalene	1.0	17
4-Chloroaniline	10	Not Detected
Hexachlorobutadiene	1.0	10
4-Chloro-3-methylphenol	5.0	Not Detected
2-Methylnaphthalene	1.0	10
Hexachlorocyclopentadiene	20	Not Detected
2,4,6-Tr chlorophenol	5.0	Not Detected
2,4,5-Tr chlorophenol	5.0	Not Detected
2-Chloronaphthalene	1.0	Not Detected
2-Nitroaniline	10	Not Detected
Dimethylphthalate	5.0	Not Detected
Acenaphthylene	1.0	Not Detected
2,6-D nitrotoluene	5.0	Not Detected
3-Nitroaniline	10	Not Detected
Acenaphthene	1.0	Not Detected
2,4-D nitrophenol	20	Not Detected
4-Nitrophenol	20	Not Detected
2,4-D nitrotoluene	5.0	Not Detected
Dibenzofuran	1.0	Not Detected
Diethylphthalate	5.0	Not Detected
Fluorene	1.0	Not Detected

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AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 3 TOX 1 INF

Lab ID#: 0612226B-03A

MODIFIED EPA METHOD TO-13A GC/MS FULL SCAN

File Name:	T12226	Date of Collection:	7/21/06
Oil Factor:	1.00	Date of Analysis:	7/21/06 10:14 PM
		Date of Extraction:	7/21/06

Compound	Rpt. Limit (μ g)	Amount (μ g)
4-Chlorophenyl-phenyl Ether	1.0	Not Detected
4-Nitroaniline	10	Not Detected
4,6-Dinitro-2-methylphenol	10	Not Detected
N-Nitrosodiphenylamine	10	Not Detected
<u>4-Bromophenyl-phenyl Ether</u>	<u>1.0</u>	<u>Not Detected</u>
Hexachlorobenzene	1.0	Not Detected
Pentachlorophenol	20	Not Detected
Phenanthrene	1.0	Not Detected
Anthracene	1.0	Not Detected
di-n-Butylphthalate	5.0	Not Detected
Fluoranthene	1.0	Not Detected
Pyrene	1.0	Not Detected
Butylbenzylphthalate	5.0	Not Detected
3,3'-Dichlorobenzidine	20	Not Detected
Chrysene	1.0	Not Detected
Benz(a)anthracene	1.0	Not Detected
bis(2-Ethylhexyl)phthalate	5.0	Not Detected
Di-n-Octylphthalate	5.0	Not Detected
Benzo(b)fluoranthene	1.0	Not Detected
Benzo(k)fluoranthene	1.0	Not Detected
Benzo(a)pyrene	1.0	Not Detected
Indeno(1,2,3-c,d)pyrene	1.0	Not Detected
Dibenz(a,h)anthracene	1.0	Not Detected
Benzo(g,h,i)perylene	1.0	Not Detected

Q = Exceeds Quality Control limits.

Container Type: XAD Tube

Surrogates	%Recovery	Method Limits
2-Fluorophenol	39 Q	50-150
Phenol-d5	96	50-150
Nitrobenzene-d5	92	50-150
2,4,6-Tribromophenol	80	50-150
Fluorene-d10	88	60-120
Pyrene-d10	98	60-120

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AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 4 TOX 1 INF DUP

Lab ID#: 0612226B-04A

MODIFIED EPA METHOD TO-13A GC/MS FULL SCAN

File Name:	06121414	Date of Collection:	12/14/06
Dil. Factor:	1.00	Date of Analysis:	12/14/06 04:44 PM
		Date of Extraction:	12/14/06

Compound	Rpt. Limit (ug)	Amount (ug)
Phenol	5.0	Not Detected
bis(2-Chloroethyl) Ether	1.0	2.7
2-Chlorophenol	5.0	Not Detected
1,3-Dichlorobenzene	1.0	3.0
1,4-Dichlorobenzene	1.0	6.0
1,2-Dichlorobenzene	1.0	25
2-Methylphenol (o-Cresol)	5.0	Not Detected
N-Nitroso-di-n-propylamine	1.0	Not Detected
4-Methylphenol/3-Methylphenol	5.0	Not Detected
Hexachloroethane	1.0	Not Detected
Nitrobenzene	1.0	Not Detected
Isophorone	1.0	2.4
2-Nitrophenol	5.0	Not Detected
2,4-Dimethylphenol	5.0	Not Detected
bis(2-Chloroethoxy) Methane	1.0	Not Detected
2,4-Dichlorophenol	5.0	Not Detected
1,2,4-Trichlorobenzene	1.0	15
Naphthalene	1.0	19
4-Chloroaniline	10	Not Detected
Hexachlorobutadiene	1.0	11
4-Chloro-3-methylphenol	5.0	Not Detected
2-Methylnaphthalene	1.0	11
Hexachlorocyclopentadiene	20	Not Detected
2,4,6-Trichlorophenol	5.0	Not Detected
2,4,5-Trichlorophenol	5.0	Not Detected
2-Chloronaphthalene	1.0	Not Detected
2-Nitroaniline	10	Not Detected
Dimethylphthalate	5.0	Not Detected
Acenaphthylene	1.0	Not Detected
2,6-Dinitrotoluene	5.0	Not Detected
3-Nitroaniline	10	Not Detected
Acenaphthene	1.0	Not Detected
2,4-Dinitrophenol	20	Not Detected
4-Nitrophenol	20	Not Detected
2,4-Dinitrotoluene	5.0	Not Detected
Dibenzofuran	1.0	Not Detected
Diethylphthalate	5.0	Not Detected
Fluorene	1.0	Not Detected



AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 4 TOX 1 INF DUP

Lab ID#: 0612226B-04A

MODIFIED EPA METHOD TO-13A GC/MS FULL SCAN

File Name:	0612226B-04A	Date of Collection: 12/14/06
Chrom Date:	12/14/06	Date of Analysis: 12/14/06 04:44:59
		Date of Extraction: 12/14/06

Compound	Rpt. Limit (ug)	Amount (ug)
4-Chlorophenyl-phenyl Ether	1.0	Not Detected
4-Nitroaniline	10	Not Detected
4,6-Dinitro-2-methylphenol	10	Not Detected
N-Nitrosodiphenylamine	10	Not Detected
4-Bromophenyl-phenyl Ether	1.0	Not Detected
Hexachlorobenzene	1.0	Not Detected
Pentachlorophenol	20	Not Detected
Phenanthrene	1.0	Not Detected
Anthracene	1.0	Not Detected
di-n-Butylphthalate	5.0	Not Detected
Fluoranthene	1.0	Not Detected
Pyrene	1.0	Not Detected
Butylbenzylphthalate	5.0	Not Detected
3,3'-Dichlorobenzidine	20	Not Detected
Chrysene	1.0	Not Detected
Benz(a)anthracene	1.0	Not Detected
bis(2-Ethylhexyl)phthalate	5.0	Not Detected
Di-n-Octylphthalate	5.0	Not Detected
Benz(b)fluoranthene	1.0	Not Detected
Benz(k)fluoranthene	1.0	Not Detected
Benzo(a)pyrene	1.0	Not Detected
Indeno(1,2,3-c,d)pyrene	1.0	Not Detected
Dibenz(a,h)anthracene	1.0	Not Detected
Benzo(g,h,i)perylene	1.0	Not Detected

Q = Exceeds Quality Control limits.

Container Type: XAD Tube

Surrogates	%Recovery	Method Limits
2-Fluorophenol	43 Q	50-150
Phenol-d5	102	50-150
Nitrobenzene-d5	94	50-150
2,4,6-Tribromophenol	90	50-150
Fluorene-d10	95	60-120
Pyrene-d10	102	60-120

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AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 5 TOX 1 EFF

Lab ID#: 0612226B-05A

MODIFIED EPA METHOD TO-13A GC/MS FULL SCAN

File Name:	P141445	Date of Collection:	12/14/06
Diff Factor:	1.00	Date of Analysis:	12/14/06 05:14 PM
		Date of Extraction:	12/14/06
Compound	Rpt. Limit (μ g)	Amount (μ g)	
Phenol	5.0		Not Detected
bis(2-Chloroethyl) Ether	1.0		Not Detected
2-Chlorophenol	5.0		Not Detected
1,3-Dichlorobenzene	1.0		Not Detected
1,4-Dichlorobenzene	1.0		Not Detected
1,2-Dichlorobenzene	1.0		Not Detected
2-Methylphenol (o-Cresol)	5.0		Not Detected
N-Nitroso-di-n-propylamine	1.0		Not Detected
4-Methylphenol/3-Methylphenol	5.0		Not Detected
Hexachloroethane	1.0		Not Detected
Nitrobenzene	1.0		Not Detected
Isophorone	1.0		Not Detected
2-Nitrophenol	5.0		Not Detected
2,4-Dimethylphenol	5.0		Not Detected
bis(2-Chloroethoxy) Methane	1.0		Not Detected
2,4-Dichlorophenol	5.0		Not Detected
1,2,4-Trichlorobenzene	1.0		Not Detected
Naphthalene	1.0		Not Detected
4-Chloroaniline	10		Not Detected
Hexachlorobutadiene	1.0		Not Detected
4-Chloro-3-methylphenol	5.0		Not Detected
2-Methylnaphthalene	1.0		Not Detected
Hexachlorocyclopentadiene	20		Not Detected
2,4,6-Trichlorophenol	5.0		Not Detected
2,4,5-Trichlorophenol	5.0		Not Detected
2-Chloronaphthalene	1.0		Not Detected
2-Nitroaniline	10		Not Detected
Dimethylphthalate	5.0		Not Detected
Acenaphthylene	1.0		Not Detected
2,6-Dinitrotoluene	5.0		Not Detected
3-Nitroaniline	10		Not Detected
Acenaphthene	1.0		Not Detected
2,4-Dinitrophenol	20		Not Detected
4-Nitrophenol	20		Not Detected
2,4-Dinitrotoluene	5.0		Not Detected
Dibenzofuran	1.0		Not Detected
Diethylphthalate	5.0		Not Detected
Fluorene	1.0		Not Detected

OKS
2/16/07



AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 5 TOX 1 EFF

Lab ID#: 0612226B-05A

MODIFIED EPA METHOD TO-13A GC/MS FULL SCAN

File Name:	5TOX1EFF	Date of Collection:	12/11/06
DL Factor:	4.00	Date of Analysis:	12/14/06 05:14 PM
		Date of Extraction:	12/12/06

Compound	Rpt. Limit (ug)	Amount (ug)
4-Chlorophenyl-phenyl Ether	1.0	Not Detected
4-Nitroaniline	10	Not Detected
4,6-Dinitro-2-methylphenol	10	Not Detected
N-Nitrosodiphenylamine	10	Not Detected
4-Bromophenyl-phenyl Ether	1.0	Not Detected
Hexachlorobenzene	1.0	Not Detected
Pentachlorophenol	20	Not Detected
Phenanthrene	1.0	Not Detected
Anthracene	1.0	Not Detected
di-n-Butylphthalate	5.0	Not Detected
Fluoranthene	1.0	Not Detected
Pyrene	1.0	Not Detected
Butylbenzylphthalate	5.0	Not Detected
3,3'-Dichlorobenzidine	20	Not Detected
Chrysene	1.0	Not Detected
Benz(a)anthracene	1.0	Not Detected
bis(2-Ethylhexyl)phthalate	5.0	Not Detected
Di-n-Octylphthalate	5.0	Not Detected
Benz(b)fluoranthene	1.0	Not Detected
Benz(k)fluoranthene	1.0	Not Detected
Benz(a)pyrene	1.0	Not Detected
Indeno(1,2,3-c,d)pyrene	1.0	Not Detected
Dibenz(a,h)anthracene	1.0	Not Detected
Benz(g,h,i)perylene	1.0	Not Detected

Container Type: XAD Tube

Surrogates	%Recovery	Method Limits
2-Fluorophenol	75	50-150
Phenol-d5	84	50-150
Nitrobenzene-d5	79	50-150
2,4,6-Tribromophenol	78	50-150
Fluorene-d10	82	60-120
Pyrene-d10	85	60-120



AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 6 TOX 2 INF

Lab ID#: 0612226B-06A

MODIFIED EPA METHOD TO-13A CC/MS FULL SCAN

Sample Name	0612226B-06A	Date of Collection	12/14/06
Dil Factor	1.00	Date of Analysis	12/14/06 08:45 PM
		Date of Extraction	12/14/06
Compound	Rpt. Limit (μ g)	Amount (μ g)	
Pheno:	5.0	Not Detected	
bis(2-Chloroethyl) Ether	1.0	3.1	
2-Chlorophenol	5.0	Not Detected	
1,3-Dichlorobenzene	1.0	1.6	
1,4-Dichlorobenzene	1.0	5.1	
1,2-Dichlorobenzene	1.0	39	
2-Methylphenol (o-Cresol)	5.0	Not Detected	
N-Nitroso-di-n-propylamine	1.0	Not Detected	
4-Methylphenol/3-Methylphenol	5.0	Not Detected	
Hexachloroethane	1.0	Not Detected	
Nitrobenzene	1.0	2.7	
Isophorone	1.0	11	
2-Nitrophenol	5.0	Not Detected	
2,4-Dimethylphenol	5.0	Not Detected	
bis(2-Chloroethoxy) Methane	1.0	Not Detected	
2,4-Dichlorophenol	5.0	Not Detected	
1,2,4-Trichlorobenzene	1.0	1.1	
Naphthalene	1.0	22	
4-Chloroaniline	10	Not Detected	
Hexachlorobutadiene	1.0	2.1	
4-Chloro-3-methylphenol	5.0	Not Detected	
2-Methylnaphthalene	1.0	4.1	
Hexachlorocyclopentadiene	20	Not Detected	
2,4,6-Trichlorophenol	5.0	Not Detected	
2,4,5-Trichlorophenol	5.0	Not Detected	
2-Chloronaphthalene	1.0	Not Detected	
2-Nitroaniline	10	Not Detected	
Dimethylphthalate	5.0	Not Detected	
Acenaphthylene	1.0	Not Detected	
2,6-Dinitrotoluene	5.0	Not Detected	
3-Nitroaniline	10	Not Detected	
Acenaphthene	1.0	Not Detected	
2,4-Dinitrophenol	20	Not Detected	
4-Nitrophenol	20	Not Detected	
2,4-Dinitrotoluene	5.0	Not Detected	
Dibenzofuran	1.0	Not Detected	
Diethylphthalate	5.0	Not Detected	
Fluorene	1.0	Not Detected	

OK
2/10



AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 6 TOX 2 INF

Lab ID#: 0612226B-06A

MODIFIED EPA METHOD TO-13A GC/MS FULL SCAN

File Name:	0612226B-06A	Date of Collection:	12/14/06
Detector:	1-PD	Date of Analysis:	12/14/06 05:44 PM
		Date of Extraction:	12/14/06

Compound	Rpt. Limit (μ g)	Amount (μ g)
4-Chlorophenyl-phenyl Ether	1.0	Not Detected
4-Nitroaniline	10	Not Detected
4,6-Dinitro-2-methylphenol	10	Not Detected
N-Nitrosodiphenylamine	10	Not Detected
<u>4-Bromophenyl-phenyl Ether</u>	1.0	Not Detected
Hexachlorobenzene	1.0	Not Detected
Pentachlorophenol	20	Not Detected
Phenanthrene	1.0	Not Detected
Anthracene	1.0	Not Detected
di-n-Butylphthalate	5.0	Not Detected
Fluoranthene	1.0	Not Detected
Pyrene	1.0	Not Detected
Butylbenzylphthalate	5.0	Not Detected
3,3'-Dichlorobenzidine	20	Not Detected
Chrysene	1.0	Not Detected
Benzo(a)anthracene	1.0	Not Detected
bis(2-Ethylhexyl)phthalate	5.0	1.3 J
Di-n-Octylphthalate	5.0	Not Detected
Benzo(b)fluoranthene	1.0	Not Detected
Benzo(k)fluoranthene	1.0	Not Detected
Benzo(a)pyrene	1.0	Not Detected
Indeno(1,2,3-c,d)pyrene	1.0	Not Detected
Dibenzo(a,h)anthracene	1.0	Not Detected
Benzo(g,h,i)perylene	1.0	Not Detected

J = Estimated value.

Q = Exceeds Quality Control limits.

Container Type: XAD Tube

Surrogates	%Recovery	Method Limits
2-Fluorophenol	13 Q	50-150
Phenol-c5	109	50-150
Nitrobenzene-d5	93	50-150
2,4,6-Tr bromophenol	88	50-150
Fluorene-d10	92	60-120
Pyrene-c10	102	60-120

OK
2/16/07



AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 7 TOX 2 INF DUP

Lab ID#: 0612226B-07A

MODIFIED EPA METHOD TO-13A GC/MS FULL SCAN

File Name:	0612226B-07A	Date of Collection:	12/11/06
File Date:	1.00	Date of Analysis:	12/14/06 14:00
		Date of Extraction:	12/11/06

Compound	Rpt. Limit (ug)	Amount (ug)
Phenol	5.0	3.2 J P
bis(2-Chloroethyl) Ether	1.0	6.1
2-Chlorophenol	5.0	Not Detected
1,3-Dichlorobenzene	1.0	3.0
1,4-Dichlorobenzene	1.0	9.1
1,2-Dichlorobenzene	1.0	72
2-Methylphenol (o-Cresol)	5.0	Not Detected
N-Nitroso-di-n-propylamine	1.0	Not Detected
4-Methylphenol/3-Methylphenol	5.0	Not Detected
Hexachloroethane	1.0	Not Detected
Nitrobenzene	1.0	Not Detected
Isophorone	1.0	22
2-Nitrophenol	5.0	Not Detected
2,4-Dimethylphenol	5.0	Not Detected
bis(2-Chloroethoxy) Methane	1.0	Not Detected
2,4-Dichlorophenol	5.0	Not Detected
1,2,4-Trichlorobenzene	1.0	2.4
Naphthalene	1.0	44
4-Chloroaniline	10	Not Detected
Hexachlorobutadiene	1.0	4.4
4-Chloro-3-methylphenol	5.0	Not Detected
2-Methylnaphthalene	1.0	8.0
Hexachlorocyclopentadiene	20	Not Detected
2,4,6-Trichlorophenol	5.0	Not Detected
2,4,5-Trichlorophenol	5.0	Not Detected
2-Chloronaphthalene	1.0	Not Detected
2-Nitroaniline	10	Not Detected
Dimethylphthalate	5.0	Not Detected
Acenaphthylene	1.0	Not Detected
2,6-Dinitrotoluene	5.0	Not Detected
3-Nitroaniline	10	Not Detected
Acenaphthene	1.0	Not Detected
2,4-Dinitrophenol	20	Not Detected
4-Nitrophenol	20	Not Detected
2,4-Dinitrotoluene	5.0	Not Detected
Dibenzofuran	1.0	Not Detected
Diethylphthalate	5.0	Not Detected
Fluorene	1.0	Not Detected



AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 7 TOX 2 INF DUP

Lab ID#: 0612226B-07A

MODIFIED EPA METHOD TO-13A GC/MS FULL SCAN

File Name:	0612226B-07A	Date of Collection:	12/11/06
Dil. Factor:	1.00	Date of Analysis:	12/12/06 15:14:39
		Date of Extraction:	12/11/06

Compound	Rpt. Limit (μ g)	Amount (μ g)
4-Chlorophenyl-phenyl Ether	1.0	Not Detected
4-Nitroaniline	10	Not Detected
4,6-Dinitro-2-methylphenol	10	Not Detected
N-Nitrosodiphenylamine	10	Not Detected
4-Bromophenyl-phenyl Ether	1.0	Not Detected
Hexachlorobenzene	1.0	Not Detected
Pentachlorophenol	20	Not Detected
Phenanthrene	1.0	Not Detected
Anthracene	1.0	Not Detected
di-n-Butylphthalate	5.0	Not Detected
Fluoranthene	1.0	Not Detected
Pyrene	1.0	Not Detected
Butylbenzylphthalate	5.0	Not Detected
3,3'-D chlorobenzidine	20	Not Detected
Chrysene	1.0	Not Detected
Benzo(a)anthracene	1.0	Not Detected
bis(2-Ethylhexyl)phthalate	5.0	1.4 J 15
Di-n-Octylphthalate	5.0	Not Detected
Benzo(b)fluoranthene	1.0	Not Detected
Benzo(k)fluoranthene	1.0	Not Detected
Benzo(a)pyrene	1.0	Not Detected
Indeno(1,2,3-c,d)pyrene	1.0	Not Detected
Dibenzo(a,h)anthracene	1.0	Not Detected
Benzo(g,h,i)perylene	1.0	Not Detected

J = Estimated value.

Q = Exceeds Quality Control limits.

Container Type: XAD Tube

Surrogates	%Recovery	Method Limits
2-Fluorophenol	17 Q	50-150
Phenol-d5	114	50-150
Nitrobenzene-d5	107	50-150
2,4,6-Tribromophenol	33	50-150
Fluorene-d10	91	60-120
Pyrene-d10	101	60-120

CH₃
26/10/06



AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 8 TOX 2 EFF

Lab ID#: 0612226B-08A

MODIFIED EPA METHOD TO-13A GC/MS FULL SCAN

File Number:	06121948	Date of Collection:	12/19/06
Div Factor:	1.00	Date of Analysis:	12/14/06 06:44 PM
		Date of Extraction:	12/12/06

Compound	Rpt. Limit (ug)	Amount (ug)
Pheno	5.0	Not Detected
bis(2-Chloroethyl) Ether	1.0	Not Detected
2-Chlorophenol	5.0	Not Detected
1,3-Dichlorobenzene	1.0	Not Detected
1,4-Dichlorobenzene	1.0	Not Detected
1,2-Dichlorobenzene	1.0	3.2
2-Methylphenol (c-Cresol)	5.0	Not Detected
N-Nitroso-di-n-propylamine	1.0	Not Detected
4-Methylphenol/3-Methylphenol	5.0	Not Detected
Hexachloroethane	1.0	Not Detected
Nitrobenzene	1.0	Not Detected
Isophorone	1.0	Not Detected
2-Nitrophenol	5.0	Not Detected
2,4-Dimethylphenol	5.0	Not Detected
bis(2-Chloroethoxy) Methane	1.0	Not Detected
2,4-Dichlorophenol	5.0	Not Detected
1,2,4-Trichlorobenzene	1.0	Not Detected
Naphthalene	1.0	4.6
4-Chloroaniline	10	Not Detected
Hexachlorobutadiene	1.0	Not Detected
4-Chloro-3-methylphenol	5.0	Not Detected
2-Methylnaphthalene	1.0	0.65 J 15
Hexachlorocyclopentadiene	20	Not Detected
2,4,6-Trichlorophenol	5.0	Not Detected
2,4,5-Trichlorophenol	5.0	Not Detected
2-Chloronaphthalene	1.0	Not Detected
2-Nitroaniline	10	Not Detected
Dimethylphthalate	5.0	Not Detected
Acenaphthylene	1.0	Not Detected
2,6-Dinitrotoluene	5.0	Not Detected
3-Nitroaniline	1.0	Not Detected
Acenaphthene	1.0	Not Detected
2,4-Dinitrophenol	20	Not Detected
4-Nitrophenol	20	Not Detected
2,4-Dinitrotoluene	5.0	Not Detected
Dibenzofuran	1.0	Not Detected
Diethyl phthalate	5.0	Not Detected
Fluorene	1.0	Not Detected

CH
JUL 16 2007
ZL



AN ENVIRONMENTAL ANALYTICAL LABORATORY

Client Sample ID: 8 TOX 2 EFF

Lab ID#: 0612226B-08A

MODIFIED EPA METHOD TO-13A GC/MS FULL SCAN

File Name:	D12226B-08A	Date of Collection:	12/11/08
Dir Factor:	1.00	Date of Analysis:	12/14/08 09:46 PM
		Date of Extraction:	12/12/08

Compound	Rpt. Limit (μ g)	Amount (μ g)
4-Chlorophenyl-phenyl Ether	1.0	Not Detected
4-Nitroaniline	10	Not Detected
4,6-Dinitro-2-methylphenol	10	Not Detected
N-Nitrosodiphenylamine	10	Not Detected
4-Bromophenyl-phenyl Ether	1.0	Not Detected
Hexachlorobenzene	1.0	Not Detected
Pentachlorophenol	20	Not Detected
Phenanthrene	1.0	Not Detected
Anthracene	1.0	Not Detected
di-n-Butylphthalate	5.0	Not Detected
Fluoranthene	1.0	Not Detected
Pyrene	1.0	Not Detected
Butylbenzylphthalate	5.0	Not Detected
3,3'-Dichlorobenzidine	20	Not Detected
Chrysene	1.0	Not Detected
Benzo(a)anthracene	1.0	Not Detected
bis(2-Ethylhexyl)phthalate	5.0	7.8
Di-n-Octylphthalate	5.0	Not Detected
Benzo(b)fluoranthene	1.0	Not Detected
Benzo(k)fluoranthene	1.0	Not Detected
Benzo(a)pyrene	1.0	Not Detected
Indeno(1,2,3-c,d)pyrene	1.0	Not Detected
Dibenz(a,h)anthracene	1.0	Not Detected
Benz(c,g,h,i)perylene	1.0	Not Detected

J = Estimated value.

Container Type: XAD Tube

Surrogates	%Recovery	Method Limits
2-Fluorophenol	84	50-150
Phenol-d5	98	50-150
Nitrobenzene-d5	91	50-150
2,4,6-Tribromophenol	88	50-150
Fluorene-d10	88	60-120
Pyrrene-d10	98	60-120

2005
2/11/09



CHAIN-OF-CUSTODY RECORD

Contact Person Chris Daily
 Company MWH Email _____
 Address 175 W. Jackson City CHICAGO State IL Zip 60601
 Phone 312 831 3415 Fax 312 831 3021

Collected by: (Signature) Chris Daily

Sample Transportation Notice

Relinquishing signature on this document indicates that sample is being shipped in compliance with all applicable local, State, Federal, national, and International laws, regulations and ordinances of any kind. Air Toxics Limited assumes no liability with respect to the collection, handling or shipping of these samples. Relinquishing signature also indicates agreement to hold harmless, defend, and indemnify Air Toxics Limited against any claim, demand, or action, of any kind, related to the collection, handling, or shipping of samples. D.O.T. Hotline (800) 467-4922.

180 BLUE RAVINE ROAD, SUITE B
 FOLSOM, CA 95630-4719
 (916) 985-1000 FAX (916) 985-1020

Page 1 of 1

Project Info:	Turn Around Time:	Lab Use Only
P.O. #	<input type="checkbox"/> Normal	Pressurized by: <u>N₂</u>
Project #	<input type="checkbox"/> Rush	Date: <u>12/11/06</u>
Project Name <u>AC5</u>	Pressurization Gas: <u>N₂</u> He	
specify		

Lab I.D.	Field Sample I.D. (Location)	Can#	Date	Time	Analyses Requested	Canister Pressure/Vacuum			
						Initial	Final	Receipt	Final (ps)
01A	1 OFFSITE ISVE	12705	12/11/06	1320	T013/T014 SIMA CAN + SORO TUBE	-30	0	0.01%	5.00
02A	2 SBPA ISVE	33702		1321		-30	0	0.01%	
03A	3 TOX 1 INF	33797		1335		-30	0	0.01%	
04A	4 TOX 1 INF DUP	29028		1345		-30	0	0.01%	
05A	5 TOX 1 EFF	911		1400		-30	0	0.01%	
06A	6 TOX 2 INF	34323		1430		-30	0	0.01%	
07A	7 TOX 2 Inf DUP	34010		1450	↓	-30	0	0.01%	
08A	8 TOX 2 EFF	31144	12-11-06	1425	T013/T014 SIMA CAN + SORO TUBE	-30	0	0.01%	5.00
						-30	0	0.01%	

Relinquished by: (signature) Date/Time <u>Chris Daily</u> 12-11-06 1500	Received by: (signature) Date/Time <u>Chris Daily</u> 12-11-06 0900	Notes:
Relinquished by: (signature) Date/Time	Received by: (signature) Date/Time	
Relinquished by: (signature) Date/Time	Received by: (signature) Date/Time	

Lab Use Only	Shipper Name	Air Bill #	Temp (°C)	Condition	Customer Seals Intact?	Work Order #
	Fed EX	8550 4759 0899	WA	good	Yes No <u>None</u>	06122264

APPENDIX C

ANNUAL SEDIMENT SAMPLE ANALYTICAL DATA

December 11, 2006

1D
GC EXTRACTABLE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

TPSS121106

Lab Name: COMPUCHEM

Contract: 8081A-8082

Lab Code: LIBRTY Case No.:

SAS No.:

SDG No.: 11759

Matrix: (soil/water) SOIL

Lab Sample ID: 1175901

Sample wt/vol: 15.0 (g/mL) G

Lab File ID: _____

% Moisture: 66 decanted: (Y/N) N

Date Received: 12/12/06

Extraction: (SepF/Cont/Sonc) SONC

Date Extracted: 12/18/06

Concentrated Extract Volume: 2500 (uL)

Date Analyzed: 12/22/06

Injection Volume: 1.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) N pH: _____

Sulfur Cleanup: (Y/N) Y

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg)	UG/KG	Q
---------	----------	---	-------	---

309-00-2-----	Aldrin	5.0	U	
319-85-7-----	beta-BHC	3.7	P	
319-84-6-----	alpha-BHC	5.0	U	
319-86-8-----	delta-BHC	5.0	U	
58-89-9-----	gamma-BHC (Lindane)	5.0	U	
72-54-8-----	4,4'-DDD	9.7	U	
72-55-9-----	4,4'-DDE	9.7	U	
50-29-3-----	4,4'-DDT	11	P	
60-57-1-----	Dieldrin	9.7	U	
959-98-8-----	Endosulfan I	5.0	U	
33213-65-9-----	Endosulfan II	9.7	U	
1031-07-8-----	Endosulfan sulfate	4.8	JP	
72-20-8-----	Endrin	9.7	U	
7421-93-4-----	Endrin Aldehyde	9.7	U	
76-44-8-----	Heptachlor	5.0	U	
1024-57-3-----	Heptachlor Epoxide	5.0	U	
72-43-5-----	Methoxychlor	50	U	
8001-35-2-----	Toxaphene	240	U	
12674-11-2-----	Aroclor-1016	71	U	
11104-28-2-----	Aroclor-1221	100	U	
11141-16-5-----	Aroclor-1232	50	U	
53469-21-9-----	Aroclor-1242	50	U	
12672-29-6-----	Aroclor-1248	300		
11097-69-1-----	Aroclor-1254	50	U	
11096-82-5-----	Aroclor-1260	50	U	
53494-70-5-----	Endrin Ketone	9.7	U	
5103-74-2-----	gamma-Chlordane	5.0	U	
5103-71-9-----	alpha-Chlordane	5.0	U	

FORM I PEST

1D
GC EXTRACTABLE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

Lab Name: COMPUCHEM

Contract: 8081A-8082

TPSSDUP

Lab Code: LIBRTY Case No.:

SAS No.: SDG No.: 11759

Matrix: (soil/water) SOIL

Lab Sample ID: 1175902

Sample wt/vol: 30.0 (g/mL) G

Lab File ID: _____

% Moisture: 55 decanted: (Y/N) N

Date Received: 12/12/06

Extraction: (SepF/Cont/Sonc) SONC

Date Extracted: 12/18/06

Concentrated Extract Volume: 5000 (uL)

Date Analyzed: 12/22/06

Injection Volume: 1.0 (uL)

Dilution Factor: 1.0

GPC Cleanup: (Y/N) N pH: _____

Sulfur Cleanup: (Y/N) Y

CAST-NOT	COMPOUND	CONCENTRATION UNITS:	
		(ug/L or ug/Kg)	UG/KG

309-00-2-----	Aldrin	2.7	JP
319-85-7-----	beta-BHC	1.9	U
319-84-6-----	alpha-BHC	3.8	U
319-86-8-----	delta-BHC	3.8	U
58-89-9-----	gamma-BHC (Lindane)	3.8	U
72-54-8-----	4,4'-DDD	7.3	U
72-55-9-----	4,4'-DDE	7.3	U
50-29-3-----	4,4'-DDT	7.3	U
50-57-1-----	Dieldrin	7.3	U
959-98-8-----	Endosulfan I	3.8	U
33213-65-9-----	Endosulfan II	6.2	JP
1031-07-8-----	Endosulfan sulfate	5.6	JP
72-20-8-----	Endrin	7.3	U
7421-93-4-----	Endrin Aldehyde	3.2	J
76-44-8-----	Heptachlor	3.8	U
1024-57-3-----	Heptachlor Epoxide	18	P
72-43-5-----	Methoxychlor	38	U
5001-35-2-----	Toxaphene	180	U
12674-11-2-----	Aroclor-1016	53	U
11104-28-2-----	Aroclor-1221	76	U
11141-16-5-----	Aroclor-1232	38	U
53469-21-9-----	Aroclor-1242	38	U
12672-29-6-----	Aroclor-1248	450	_____
11097-69-1-----	Aroclor-1254	38	U
11096-82-5-----	Aroclor-1260	38	U
53494-70-5-----	Endrin Ketone	7.3	U
5103-74-2-----	gamma-Chlordane	11	P
5103-71-9-----	alpha-Chlordane	3.8	U

FORM I PEST

11/24/07